



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 27, 2026 – 02:53 PM UTC

PDB ID : 6Q2D / pdb_00006q2d
Title : Crystal structure of Methanobrevibacter smithii Dph2 in complex with Methanobrevibacter smithii elongation factor 2
Authors : Fenwick, M.K.; Dong, M.; Lin, H.; Ealick, S.E.
Deposited on : 2019-08-07
Resolution : 3.45 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

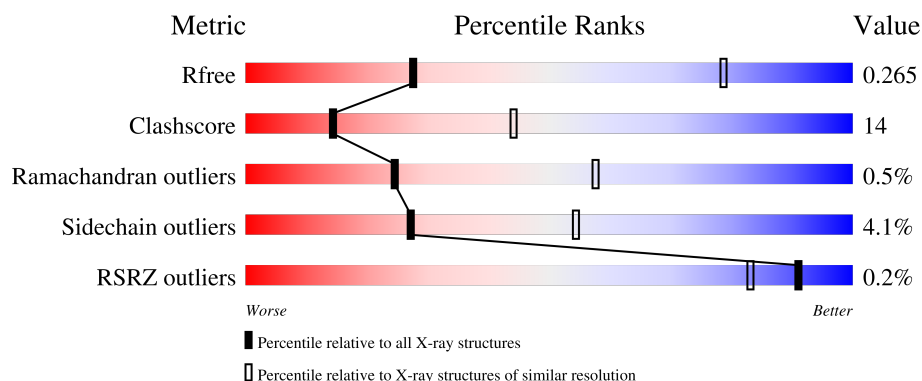
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.45 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))
R_{free}	180053	1070 (3.50-3.42)
Clashscore	190562	1128 (3.50-3.42)
Ramachandran outliers	187476	1101 (3.50-3.42)
Sidechain outliers	187428	1102 (3.50-3.42)
RSRZ outliers	180081	1070 (3.50-3.42)

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

Mol	Chain	Length	Quality of chain
1	A	337	% 77% 19% ..
1	B	337	68% 26% ..
2	C	733	61% 31% • 5%
2	F	733	41% 56%

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
3	SF4	A	401	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12185 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 2-(3-amino-3-carboxypropyl)histidine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	0	0
			2586	1661	427	485	13			
1	B	325	Total	C	N	O	S	0	0	0
			2570	1652	423	482	13			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP A5UMY5
A	-1	SER	-	expression tag	UNP A5UMY5
A	0	HIS	-	expression tag	UNP A5UMY5
B	-2	GLY	-	expression tag	UNP A5UMY5
B	-1	SER	-	expression tag	UNP A5UMY5
B	0	HIS	-	expression tag	UNP A5UMY5

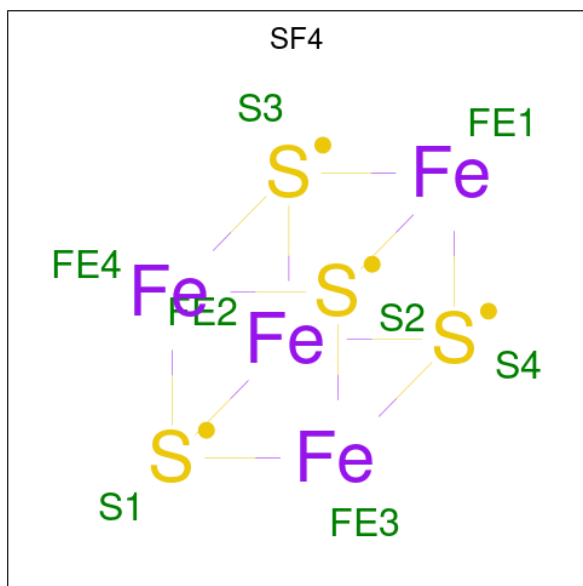
- Molecule 2 is a protein called Elongation factor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	700	Total	C	N	O	S	0	0	0
			5406	3408	922	1046	30			
2	F	326	Total	C	N	O		0	0	0
			1607	955	326	326				

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	expression tag	UNP A0A2H4U7K7
C	-1	SER	-	expression tag	UNP A0A2H4U7K7
C	0	GLY	-	expression tag	UNP A0A2H4U7K7
F	-2	GLY	-	expression tag	UNP A0A2H4U7K7
F	-1	SER	-	expression tag	UNP A0A2H4U7K7
F	0	GLY	-	expression tag	UNP A0A2H4U7K7

- Molecule 3 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).

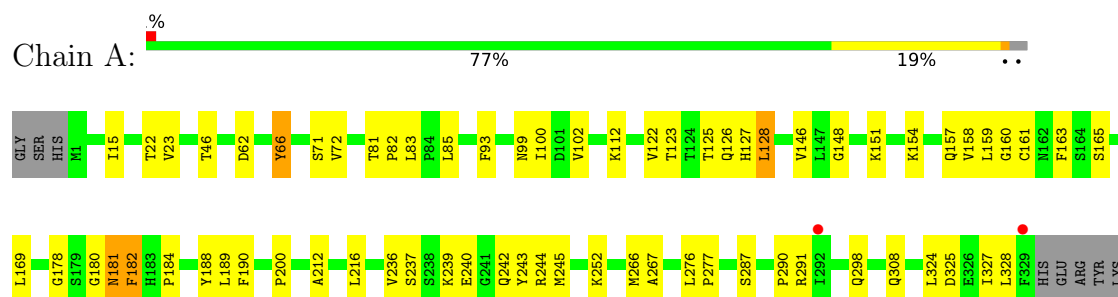


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Fe	S	0	0
			8	4	4		
3	B	1	Total	Fe	S	0	0
			8	4	4		

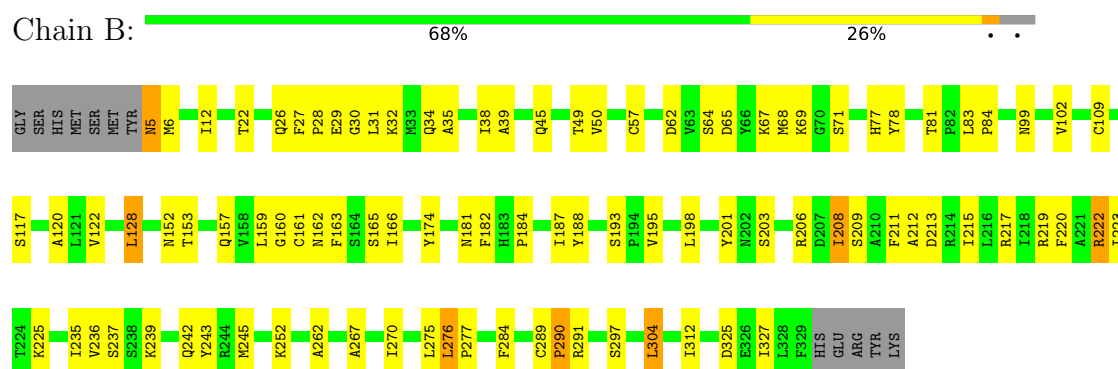
3 Residue-property plots

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 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.

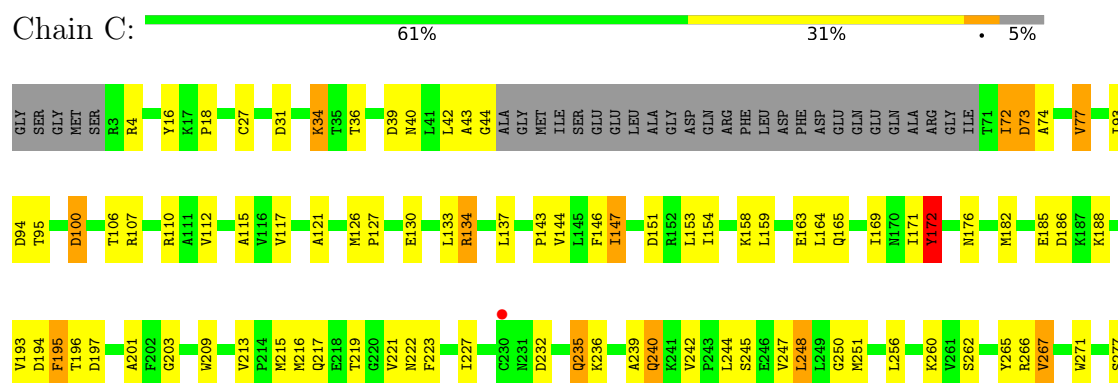
- Molecule 1: 2-(3-amino-3-carboxypropyl)histidine synthase

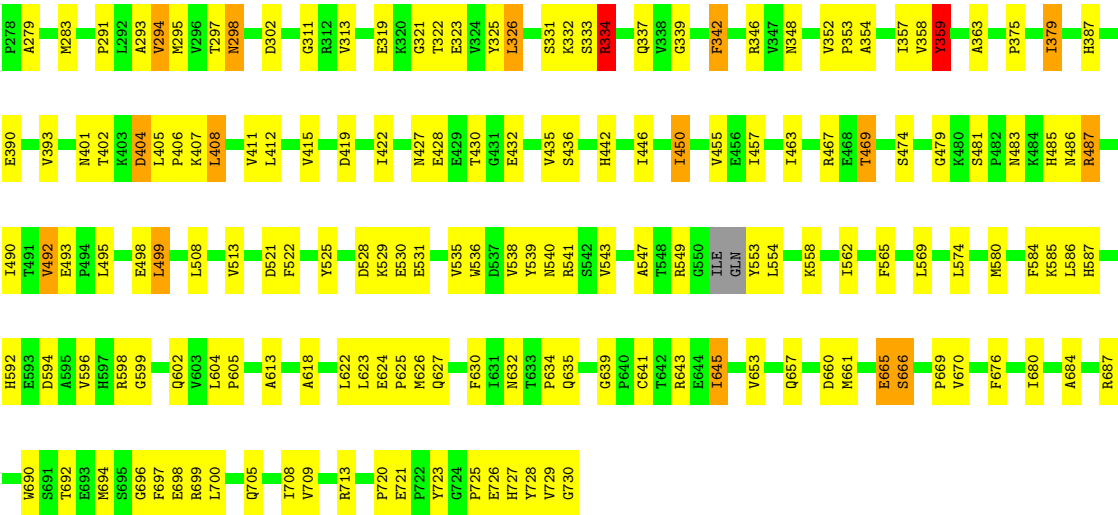


- Molecule 1: 2-(3-amino-3-carboxypropyl)histidine synthase

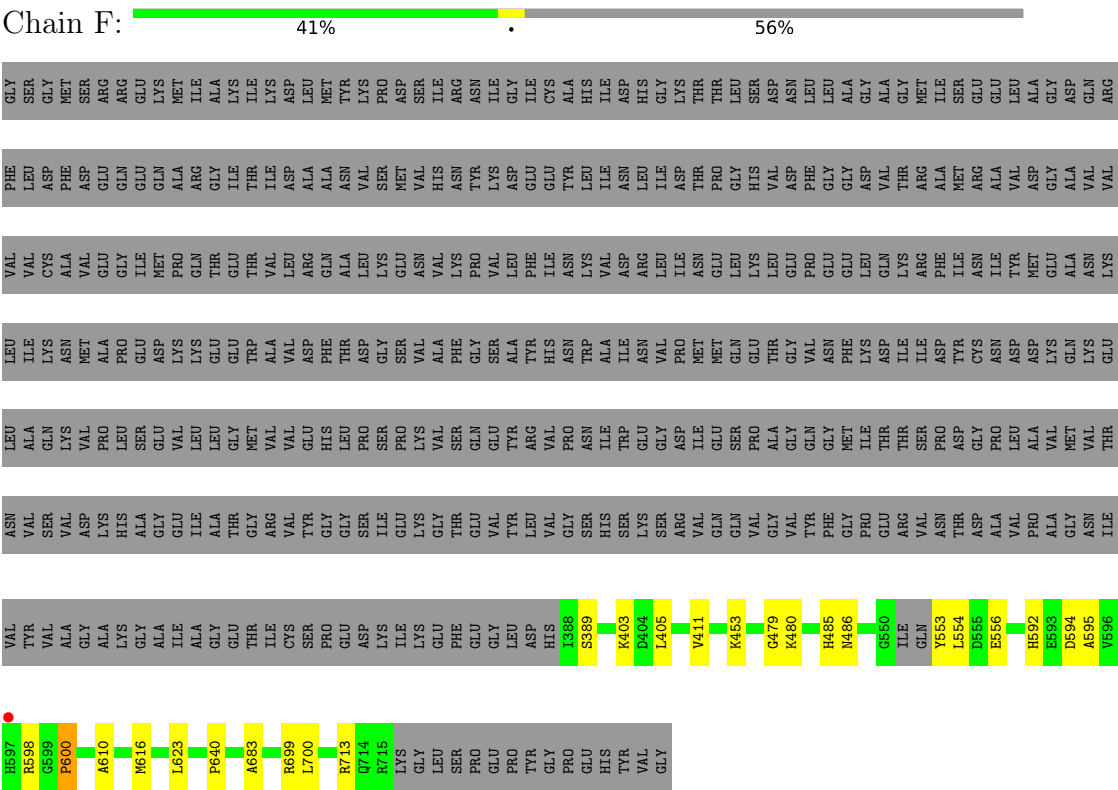


- Molecule 2: Elongation factor 2





● Molecule 2: Elongation factor 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	100.90Å 322.98Å 172.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.36 – 3.45 49.36 – 3.45	Depositor EDS
% Data completeness (in resolution range)	99.3 (49.36-3.45) 99.6 (49.36-3.45)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 3.48Å)	Xtriage
Refinement program	PHENIX 1.14_3211	Depositor
R, R_{free}	0.224 , 0.267 0.227 , 0.265	Depositor DCC
R_{free} test set	1870 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	139.3	Xtriage
Anisotropy	0.652	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 194.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12185	wwPDB-VP
Average B, all atoms (Å ²)	208.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.84% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SF4

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	0.69	0/2633	1.03	6/3561 (0.2%)
1	B	0.66	0/2617	1.00	1/3538 (0.0%)
2	C	0.85	0/5502	1.27	17/7442 (0.2%)
2	F	0.92	0/1605	1.19	0/2230
All	All	0.79	0/12357	1.15	24/16771 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	2
2	C	0	4
2	F	0	1
All	All	0	8

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	182	PHE	N-CA-CB	9.14	123.27	110.01
2	C	240	GLN	CA-C-N	-9.09	108.42	122.60
2	C	240	GLN	C-N-CA	-9.09	108.42	122.60
2	C	172	TYR	CB-CA-C	7.28	125.24	110.38
2	C	195	PHE	CA-C-N	6.74	129.64	120.54
2	C	195	PHE	C-N-CA	6.74	129.64	120.54
1	A	181	ASN	CA-C-N	-5.97	112.68	120.44
1	A	181	ASN	C-N-CA	-5.97	112.68	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	541	ARG	CB-CA-C	-5.77	103.21	112.09
2	C	599	GLY	CA-C-N	5.76	125.43	119.56
2	C	599	GLY	C-N-CA	5.76	125.43	119.56
1	B	290	PRO	CB-CA-C	-5.65	102.24	111.56
2	C	250	GLY	CA-C-N	5.52	127.68	120.28
2	C	250	GLY	C-N-CA	5.52	127.68	120.28
2	C	34	LYS	CA-C-N	5.50	128.44	120.79
2	C	34	LYS	C-N-CA	5.50	128.44	120.79
2	C	359	TYR	CB-CA-C	5.46	119.17	109.72
1	A	182	PHE	CB-CA-C	-5.43	102.36	110.88
1	A	46	THR	O-C-N	-5.26	118.34	121.71
2	C	186	ASP	CA-CB-CG	5.25	117.85	112.60
1	A	66	TYR	CB-CA-C	-5.25	102.64	110.88
2	C	334	ARG	CB-CA-C	5.15	118.70	110.16
2	C	267	VAL	O-C-N	-5.14	117.13	120.42
2	C	232	ASP	CA-CB-CG	5.04	117.64	112.60

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	178	GLY	Peptide
1	B	102	VAL	Peptide
1	B	5	ASN	Peptide
2	C	134	ARG	Sidechain
2	C	302	ASP	Peptide
2	C	4	ARG	Sidechain
2	C	486	ASN	Peptide
2	F	556	GLU	Peptide

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	2586	0	2617	59	0
1	B	2570	0	2606	72	0
2	C	5406	0	5367	196	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	1607	0	723	16	0
3	A	8	0	0	3	0
3	B	8	0	0	1	0
All	All	12185	0	11313	332	1

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 14.

All (332) 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:181:ASN:OD1	1:B:206:ARG:NH2	1.97	0.96
2:F:640:PRO:CB	2:F:683:ALA:O	2.14	0.95
1:A:161:CYS:SG	3:A:401:SF4:FE2	1.63	0.91
1:B:211:PHE:CE2	1:B:215:ILE:HD11	2.13	0.83
1:B:225:LYS:HG2	2:C:428:GLU:OE2	1.79	0.82
2:C:159:LEU:HD22	2:C:163:GLU:HB3	1.65	0.78
2:C:313:VAL:HG23	2:C:358:VAL:CG2	2.14	0.77
2:C:467:ARG:HH11	2:C:580:MET:HB3	1.48	0.76
2:C:325:TYR:HD1	2:C:331:SER:O	1.70	0.75
2:C:626:MET:SD	2:C:698:GLU:HG3	2.27	0.75
1:A:146:VAL:HG21	1:A:169:LEU:HD22	1.68	0.75
1:A:161:CYS:HG	3:A:401:SF4:FE2	1.03	0.75
2:C:134:ARG:NE	2:C:182:MET:HE3	2.01	0.74
2:C:645:ILE:HD13	2:C:666:SER:OG	1.87	0.74
2:C:40:ASN:O	2:C:43:ALA:HB3	1.87	0.73
2:C:411:VAL:HG21	2:C:455:VAL:HG23	1.72	0.72
2:C:538:VAL:HG22	2:C:543:VAL:HG12	1.71	0.71
1:A:161:CYS:SG	3:A:401:SF4:S4	2.89	0.71
2:C:279:ALA:HA	2:C:291:PRO:HG2	1.74	0.70
2:C:513:VAL:HG21	2:C:535:VAL:HG13	1.74	0.70
2:C:262:SER:HA	2:C:265:TYR:CE2	2.27	0.69
2:C:346:ARG:HD2	2:C:359:TYR:CD2	2.28	0.69
2:C:134:ARG:HE	2:C:182:MET:HE3	1.58	0.69
2:C:337:GLN:NE2	2:C:348:ASN:HB3	2.07	0.69
2:C:450:ILE:HG21	2:C:457:ILE:CD1	2.23	0.69
2:C:342:PHE:HD2	2:C:342:PHE:O	1.76	0.68
1:B:161:CYS:SG	1:B:327:ILE:HB	2.33	0.68
2:C:154:ILE:HG23	2:C:227:ILE:HG23	1.75	0.67
2:C:408:LEU:HA	2:C:455:VAL:HG21	1.77	0.67
1:A:308:GLN:OE1	1:A:325:ASP:HB2	1.94	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:LEU:HD22	1:A:216:LEU:HD21	1.78	0.66
1:A:81:THR:HG21	1:A:154:LYS:HD2	1.78	0.66
2:C:130:GLU:O	2:C:134:ARG:HG3	1.97	0.65
1:A:102:VAL:HG12	1:A:200:PRO:HB3	1.79	0.65
1:B:290:PRO:HB3	1:B:325:ASP:HB3	1.78	0.65
1:A:163:PHE:HD1	1:A:190:PHE:CD2	2.15	0.64
1:B:211:PHE:CE2	1:B:215:ILE:CD1	2.81	0.63
2:C:16:TYR:CD1	2:C:353:PRO:HG3	2.33	0.63
1:B:211:PHE:CZ	1:B:215:ILE:HD11	2.33	0.63
1:B:5:ASN:OD1	1:B:6:MET:N	2.32	0.63
1:A:160:GLY:HA2	1:A:182:PHE:HE2	1.64	0.63
1:A:244:ARG:HD3	1:A:287:SER:O	1.98	0.63
2:C:34:LYS:NZ	2:C:95:THR:O	2.32	0.63
1:A:163:PHE:CD1	1:A:190:PHE:CD2	2.87	0.62
1:A:243:TYR:CZ	1:A:245:MET:HB2	2.35	0.62
2:C:467:ARG:NE	2:C:580:MET:SD	2.71	0.62
2:F:623:LEU:HA	2:F:700:LEU:H	1.64	0.62
2:C:492:VAL:CG1	2:C:584:PHE:CE2	2.84	0.61
2:C:531:GLU:OE2	2:C:585:LYS:HE2	2.01	0.61
1:A:82:PRO:HB3	1:A:93:PHE:CZ	2.36	0.60
2:C:467:ARG:HH11	2:C:580:MET:CB	2.14	0.60
1:A:182:PHE:CB	1:A:291:ARG:HH21	2.15	0.60
2:C:553:TYR:O	2:C:554:LEU:HB2	2.02	0.60
1:A:244:ARG:HH12	1:A:325:ASP:CG	2.09	0.59
1:B:12:ILE:HD13	1:B:45:GLN:OE1	2.03	0.59
2:C:337:GLN:HG2	2:C:348:ASN:HD22	1.67	0.59
2:C:342:PHE:O	2:C:342:PHE:CD2	2.55	0.59
2:C:182:MET:HE1	2:C:661:MET:HB3	1.85	0.59
2:C:485:HIS:HB3	2:C:592:HIS:CB	2.33	0.58
1:A:123:THR:HG23	1:A:128:LEU:HD13	1.85	0.58
2:C:450:ILE:HG21	2:C:457:ILE:HD13	1.86	0.58
1:A:163:PHE:HD1	1:A:190:PHE:CG	2.21	0.58
2:C:72:ILE:HD13	2:C:77:VAL:HG21	1.85	0.58
1:A:160:GLY:CA	1:A:182:PHE:HE2	2.17	0.58
1:B:122:VAL:HG22	1:B:166:ILE:HG21	1.86	0.58
1:A:123:THR:OG1	1:A:127:HIS:HB2	2.03	0.57
2:C:337:GLN:CG	2:C:348:ASN:HD22	2.18	0.57
1:B:81:THR:HG21	1:B:157:GLN:HE22	1.69	0.57
2:C:634:PRO:HG2	2:C:687:ARG:HD3	1.85	0.57
2:C:73:ASP:N	2:C:73:ASP:OD1	2.38	0.57
2:C:339:GLY:HA3	2:C:359:TYR:CE1	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:632:ASN:HB3	2:C:661:MET:CE	2.34	0.57
2:C:279:ALA:HA	2:C:291:PRO:CG	2.34	0.57
1:A:188:TYR:CD2	1:A:212:ALA:CB	2.88	0.56
1:B:219:ARG:HH12	1:B:297:SER:CB	2.17	0.56
2:C:574:LEU:HD21	2:C:705:GLN:NE2	2.19	0.56
2:C:325:TYR:HB3	2:C:375:PRO:HA	1.86	0.56
1:B:120:ALA:HB3	1:B:174:TYR:CD1	2.41	0.56
2:C:508:LEU:HD21	2:C:521:ASP:HB3	1.87	0.56
2:C:144:VAL:HG23	2:C:256:LEU:HD21	1.88	0.55
1:A:82:PRO:HB3	1:A:93:PHE:CE2	2.42	0.55
2:C:558:LYS:O	2:C:562:ILE:HG12	2.06	0.55
2:C:613:ALA:HB2	2:C:728:TYR:CB	2.35	0.55
1:B:160:GLY:O	1:B:182:PHE:HE2	1.90	0.55
2:C:159:LEU:HD12	2:C:164:LEU:HD12	1.88	0.55
2:C:339:GLY:HA3	2:C:359:TYR:HE1	1.70	0.55
1:B:239:LYS:HD2	1:B:242:GLN:NE2	2.22	0.55
2:C:334:ARG:NH2	2:C:334:ARG:HG3	2.22	0.55
2:C:159:LEU:HB2	2:C:164:LEU:HD12	1.89	0.54
1:B:212:ALA:O	1:B:213:ASP:C	2.51	0.54
1:B:213:ASP:O	1:B:217:ARG:HG3	2.08	0.54
2:C:100:ASP:HB3	2:C:690:TRP:HE1	1.73	0.54
2:C:540:ASN:O	2:C:580:MET:HA	2.07	0.54
1:B:284:PHE:HB2	1:B:304:LEU:HD12	1.88	0.54
2:C:293:ALA:HA	2:C:379:ILE:HD13	1.90	0.53
2:C:479:GLY:O	2:C:487:ARG:HA	2.08	0.53
2:C:185:GLU:OE1	2:C:188:LYS:CE	2.56	0.53
1:B:26:GLN:OE1	1:B:77:HIS:ND1	2.40	0.53
2:C:313:VAL:HG23	2:C:358:VAL:HG21	1.89	0.53
2:C:352:VAL:HG11	2:C:358:VAL:HG22	1.91	0.53
1:A:125:THR:OG1	1:A:157:GLN:NE2	2.42	0.53
2:C:401:ASN:O	2:C:404:ASP:OD2	2.26	0.53
2:C:639:GLY:O	2:C:643:ARG:HG3	2.09	0.53
1:B:222:ARG:NH2	1:B:297:SER:HB2	2.23	0.53
2:C:107:ARG:HA	2:C:110:ARG:HD3	1.90	0.53
2:C:121:ALA:HB2	2:C:147:ILE:HG22	1.90	0.53
2:C:271:TRP:CD2	2:C:283:MET:HE1	2.44	0.52
2:C:490:ILE:HG22	2:C:586:LEU:HA	1.89	0.52
2:C:547:ALA:HB3	2:C:587:HIS:HA	1.91	0.52
1:B:29:GLU:HG2	1:B:32:LYS:HE2	1.90	0.52
2:C:18:PRO:HB2	2:C:260:LYS:HB2	1.92	0.52
2:C:522:PHE:CZ	2:C:535:VAL:HG11	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:422:ILE:HD11	2:C:435:VAL:HG12	1.91	0.52
2:C:481:SER:OG	2:C:483:ASN:OD1	2.28	0.52
2:C:565:PHE:CE1	2:C:569:LEU:HD11	2.45	0.52
1:A:125:THR:HA	1:A:128:LEU:HD22	1.91	0.52
1:A:99:ASN:O	1:A:100:ILE:HD13	2.09	0.52
1:A:126:GLN:NE2	2:C:596:VAL:HG11	2.24	0.52
2:F:623:LEU:HA	2:F:700:LEU:N	2.25	0.52
1:A:182:PHE:HB2	1:A:291:ARG:HH21	1.75	0.52
1:B:276:LEU:N	1:B:277:PRO:HD2	2.25	0.51
1:B:327:ILE:HD12	3:B:401:SF4:S4	2.50	0.51
2:C:528:ASP:OD1	2:C:529:LYS:N	2.44	0.51
2:C:632:ASN:HB3	2:C:661:MET:HE2	1.92	0.51
2:C:195:PHE:CE1	2:C:213:VAL:HG23	2.45	0.51
1:A:276:LEU:HB3	2:F:600:PRO:CB	2.41	0.51
1:B:188:TYR:CE2	1:B:209:SER:HA	2.46	0.51
2:C:321:GLY:HA2	2:C:334:ARG:HH11	1.76	0.51
2:C:332:LYS:HG2	2:C:333:SER:N	2.26	0.51
1:A:188:TYR:CD2	1:A:212:ALA:HB1	2.46	0.51
1:B:38:ILE:HD13	1:B:78:TYR:OH	2.11	0.51
1:B:109:CYS:SG	1:B:198:LEU:HD22	2.51	0.51
2:C:627:GLN:HG3	2:C:670:VAL:HG22	1.93	0.50
2:C:694:MET:HE3	2:C:696:GLY:HA2	1.93	0.50
1:A:182:PHE:CE1	1:A:291:ARG:HG3	2.46	0.50
2:C:16:TYR:CE1	2:C:353:PRO:HG3	2.46	0.50
2:C:213:VAL:HA	2:C:216:MET:CE	2.41	0.50
1:A:66:TYR:HE2	1:A:240:GLU:OE1	1.93	0.50
1:B:222:ARG:HH21	1:B:297:SER:HB2	1.76	0.50
1:B:239:LYS:HD2	1:B:242:GLN:HE21	1.76	0.50
2:C:215:MET:HE3	2:C:247:VAL:HG22	1.92	0.50
2:C:93:ILE:HD11	2:C:357:ILE:HG13	1.94	0.50
2:C:154:ILE:CG2	2:C:227:ILE:HG23	2.39	0.50
2:C:720:PRO:HB2	2:C:721:GLU:OE1	2.12	0.50
2:C:236:LYS:O	2:C:240:GLN:HG2	2.12	0.50
2:C:723:TYR:HD1	2:C:727:HIS:CD2	2.30	0.50
1:B:161:CYS:SG	1:B:327:ILE:HD12	2.52	0.49
1:A:123:THR:HG21	1:A:128:LEU:HA	1.95	0.49
2:C:151:ASP:HB3	2:C:235:GLN:OE1	2.13	0.49
2:C:342:PHE:CD2	2:C:342:PHE:C	2.90	0.49
2:F:480:LYS:HA	2:F:486:ASN:O	2.13	0.49
2:C:547:ALA:HB3	2:C:586:LEU:O	2.12	0.49
1:A:252:LYS:HE2	1:B:49:THR:HG23	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:625:PRO:HB3	2:C:697:PHE:CE1	2.47	0.49
1:A:298:GLN:NE2	2:C:483:ASN:O	2.43	0.49
2:C:492:VAL:CG1	2:C:584:PHE:CD2	2.95	0.49
1:A:290:PRO:HB3	1:A:325:ASP:HB3	1.95	0.48
1:A:276:LEU:N	1:A:277:PRO:CD	2.76	0.48
1:B:217:ARG:NH1	2:C:402:THR:HG23	2.29	0.48
2:C:223:PHE:O	2:C:227:ILE:HG13	2.13	0.48
1:A:181:ASN:O	1:A:182:PHE:C	2.55	0.48
2:C:313:VAL:CG2	2:C:358:VAL:HG21	2.43	0.48
1:B:219:ARG:HH12	1:B:297:SER:HB2	1.79	0.48
2:C:219:THR:OG1	2:C:221:VAL:HG23	2.13	0.48
2:C:474:SER:HB3	2:C:618:ALA:HB2	1.95	0.48
2:C:729:VAL:HG12	2:C:730:GLY:N	2.27	0.48
2:C:326:LEU:HD21	2:C:363:ALA:O	2.14	0.48
2:C:700:LEU:HD21	2:C:708:ILE:HD12	1.95	0.48
1:B:184:PRO:HA	1:B:187:ILE:HD12	1.96	0.48
2:C:147:ILE:HD13	2:C:171:ILE:HG21	1.96	0.48
2:C:598:ARG:HA	2:C:602:GLN:OE1	2.13	0.48
1:A:182:PHE:CB	1:A:291:ARG:NH2	2.76	0.47
2:C:334:ARG:HG3	2:C:334:ARG:HH21	1.79	0.47
2:C:294:VAL:HG23	2:C:295:MET:N	2.28	0.47
2:C:635:GLN:OE1	2:C:657:GLN:NE2	2.47	0.47
2:C:44:GLY:HA3	2:C:245:SER:OG	2.14	0.47
1:B:237:SER:HA	1:B:267:ALA:O	2.14	0.47
2:C:27:CYS:SG	2:C:115:ALA:HB1	2.55	0.47
2:C:159:LEU:HB2	2:C:164:LEU:CD1	2.43	0.47
2:C:415:VAL:HG11	2:C:446:ILE:HG23	1.97	0.47
2:C:422:ILE:HD12	2:C:436:SER:O	2.13	0.47
1:A:62:ASP:OD1	1:A:239:LYS:HG2	2.15	0.47
1:A:237:SER:HA	1:A:267:ALA:O	2.13	0.47
1:B:152:ASN:HB2	1:B:162:ASN:HD21	1.79	0.47
2:C:203:GLY:HA2	2:C:209:TRP:CZ3	2.49	0.47
2:C:522:PHE:CE2	2:C:535:VAL:HG11	2.49	0.47
1:A:148:GLY:CA	1:A:169:LEU:HD11	2.45	0.47
1:B:5:ASN:OD1	1:B:5:ASN:C	2.58	0.47
2:C:313:VAL:CG1	2:C:354:ALA:HA	2.45	0.47
2:C:427:ASN:ND2	2:C:430:THR:OG1	2.48	0.46
2:C:531:GLU:OE2	2:C:585:LYS:CE	2.63	0.46
2:C:553:TYR:O	2:C:554:LEU:CB	2.63	0.46
2:F:411:VAL:CB	2:F:453:LYS:CB	2.93	0.46
2:C:222:ASN:C	2:C:222:ASN:OD1	2.59	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:298:ASN:C	2:C:298:ASN:OD1	2.57	0.46
1:B:153:THR:OG1	1:B:165:SER:HB3	2.14	0.46
1:B:181:ASN:HA	1:B:184:PRO:CG	2.46	0.46
1:B:276:LEU:N	1:B:277:PRO:CD	2.79	0.46
2:C:297:THR:HG23	2:C:311:GLY:HA2	1.96	0.46
1:A:236:VAL:O	1:A:266:MET:HA	2.16	0.46
2:C:201:ALA:HB3	2:C:251:MET:HE1	1.97	0.46
2:C:407:LYS:O	2:C:411:VAL:HG23	2.15	0.46
2:C:169:ILE:HA	2:C:172:TYR:CD2	2.51	0.46
1:A:236:VAL:HA	1:A:287:SER:HG	1.81	0.46
2:C:622:LEU:HD22	2:C:709:VAL:HG22	1.97	0.46
1:A:15:ILE:HG12	1:A:23:VAL:HG21	1.98	0.45
2:C:100:ASP:HB3	2:C:690:TRP:NE1	2.32	0.45
1:A:148:GLY:HA3	1:A:165:SER:O	2.16	0.45
1:A:239:LYS:HD2	1:A:242:GLN:CD	2.42	0.45
1:B:225:LYS:CG	2:C:428:GLU:OE2	2.56	0.45
1:B:290:PRO:HG3	1:B:325:ASP:CG	2.41	0.45
2:C:185:GLU:OE1	2:C:188:LYS:HE2	2.17	0.45
1:B:57:CYS:HA	1:B:62:ASP:OD2	2.16	0.45
1:B:223:ILE:HG21	1:B:312:ILE:HG22	1.98	0.45
2:C:154:ILE:HD13	2:C:227:ILE:HA	1.97	0.45
2:C:194:ASP:OD1	2:C:196:THR:HB	2.16	0.45
2:C:319:GLU:HB2	2:C:322:THR:OG1	2.17	0.45
2:C:326:LEU:N	2:C:326:LEU:CD1	2.79	0.45
1:B:220:PHE:HA	1:B:223:ILE:HD12	1.98	0.45
2:C:172:TYR:CE2	2:C:195:PHE:HB2	2.52	0.45
2:C:624:GLU:N	2:C:698:GLU:O	2.50	0.45
1:B:81:THR:CG2	1:B:157:GLN:HE22	2.28	0.45
2:C:492:VAL:HG11	2:C:584:PHE:CE2	2.52	0.45
2:C:334:ARG:HH21	2:C:334:ARG:CG	2.30	0.45
1:A:112:LYS:HD2	1:A:112:LYS:HA	1.70	0.45
1:B:128:LEU:HD11	1:B:157:GLN:HB2	1.99	0.45
2:C:393:VAL:HG13	2:C:463:ILE:O	2.17	0.45
2:C:676:PHE:CZ	2:C:680:ILE:HD12	2.52	0.45
1:A:188:TYR:HD2	1:A:212:ALA:HB1	1.82	0.44
1:B:83:LEU:HB3	1:B:84:PRO:HD2	1.99	0.44
2:C:74:ALA:O	2:C:94:ASP:HB2	2.17	0.44
2:C:419:ASP:O	2:C:422:ILE:HG22	2.17	0.44
2:C:690:TRP:HZ3	2:C:692:THR:HG23	1.82	0.44
2:C:721:GLU:OE2	2:F:616:MET:O	2.34	0.44
1:B:27:PHE:CE2	1:B:35:ALA:HB2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:159:LEU:HB3	2:C:163:GLU:HB2	1.98	0.44
2:C:498:GLU:HB2	2:C:525:TYR:O	2.17	0.44
1:A:146:VAL:CG2	1:A:169:LEU:HD22	2.44	0.44
2:C:405:LEU:HB3	2:C:406:PRO:HD3	2.00	0.44
1:B:22:THR:HG21	1:B:71:SER:O	2.17	0.44
2:C:126:MET:HB3	2:C:127:PRO:HD2	1.99	0.44
2:C:411:VAL:HG21	2:C:455:VAL:CG2	2.45	0.44
2:C:469:THR:HG22	2:C:623:LEU:HD11	1.99	0.44
2:C:604:LEU:N	2:C:605:PRO:HD2	2.31	0.44
2:F:403:LYS:C	2:F:405:LEU:H	2.25	0.44
2:F:485:HIS:O	2:F:592:HIS:CB	2.66	0.44
1:A:22:THR:HG22	1:A:72:VAL:HA	2.00	0.44
2:C:112:VAL:O	2:C:266:ARG:CZ	2.66	0.44
1:A:180:GLY:O	1:A:184:PRO:HD2	2.17	0.44
1:B:28:PRO:HD2	1:B:31:LEU:HD12	2.00	0.44
1:B:235:ILE:HG21	1:B:270:ILE:HG12	2.00	0.44
2:C:137:LEU:HD21	2:C:143:PRO:HG3	1.99	0.43
2:C:72:ILE:HG22	2:C:73:ASP:N	2.32	0.43
2:C:159:LEU:CD1	2:C:164:LEU:HD12	2.46	0.43
2:C:528:ASP:HB3	2:C:531:GLU:HG3	1.99	0.43
2:C:626:MET:HG2	2:C:669:PRO:HA	2.00	0.43
1:B:223:ILE:HG21	1:B:312:ILE:CG2	2.48	0.43
2:C:705:GLN:O	2:C:709:VAL:HG23	2.19	0.43
2:C:144:VAL:HG23	2:C:256:LEU:CD2	2.47	0.43
2:F:479:GLY:HA3	2:F:610:ALA:HB2	1.99	0.43
1:A:83:LEU:HB3	1:A:85:LEU:HG	2.01	0.43
1:A:122:VAL:CG1	1:A:158:VAL:HG13	2.48	0.43
1:B:289:CYS:SG	1:B:291:ARG:HB2	2.59	0.43
2:C:134:ARG:NH2	2:C:634:PRO:HG3	2.34	0.43
1:B:39:ALA:HA	1:B:50:VAL:HG21	2.00	0.43
2:C:93:ILE:HD11	2:C:357:ILE:CG1	2.49	0.43
2:C:630:PHE:O	2:C:690:TRP:HB2	2.19	0.43
1:B:208:ILE:HD12	1:B:208:ILE:HA	1.85	0.43
2:C:146:PHE:CD2	2:C:248:LEU:HD11	2.53	0.43
1:B:159:LEU:HB2	1:B:162:ASN:HB3	2.00	0.43
1:B:290:PRO:CB	1:B:325:ASP:HB3	2.46	0.43
2:C:213:VAL:HA	2:C:216:MET:HE2	2.01	0.43
1:A:159:LEU:O	1:A:163:PHE:CE2	2.72	0.42
2:C:209:TRP:HB3	2:C:239:ALA:HA	2.01	0.42
2:C:613:ALA:HB2	2:C:728:TYR:HB2	2.00	0.42
1:A:239:LYS:HB2	1:A:242:GLN:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:485:HIS:HB3	2:C:592:HIS:HB2	2.01	0.42
2:C:726:GLU:HG2	2:F:713:ARG:CB	2.49	0.42
1:B:184:PRO:HB2	1:B:208:ILE:HG12	2.00	0.42
2:C:323:GLU:HG3	2:C:334:ARG:HB2	2.02	0.42
1:B:160:GLY:O	1:B:163:PHE:HE2	2.02	0.42
2:C:93:ILE:HD11	2:C:357:ILE:HD11	1.99	0.42
2:C:134:ARG:HE	2:C:182:MET:CE	2.30	0.42
2:C:165:GLN:OE1	2:C:223:PHE:CZ	2.72	0.42
2:C:694:MET:HE3	2:C:696:GLY:CA	2.50	0.42
1:B:30:GLY:HA2	2:F:598:ARG:O	2.19	0.42
2:C:39:ASP:O	2:C:42:LEU:HB3	2.20	0.42
2:C:641:CYS:SG	2:C:684:ALA:HB2	2.59	0.42
1:A:243:TYR:CZ	1:A:245:MET:CB	3.01	0.42
1:B:161:CYS:HB3	1:B:290:PRO:HG2	2.02	0.42
1:B:252:LYS:NZ	1:B:262:ALA:O	2.53	0.42
1:A:163:PHE:CE1	1:A:324:LEU:HD12	2.55	0.42
2:C:182:MET:HE1	2:C:661:MET:CB	2.50	0.42
2:C:213:VAL:HA	2:C:216:MET:HE3	2.02	0.42
2:C:574:LEU:O	2:C:713:ARG:HD2	2.20	0.42
1:B:181:ASN:CG	1:B:206:ARG:HH21	2.20	0.42
2:C:405:LEU:HA	2:C:405:LEU:HD12	1.87	0.42
1:A:81:THR:HG21	1:A:154:LYS:CD	2.49	0.41
2:C:121:ALA:HB1	2:C:153:LEU:HD11	2.02	0.41
2:C:635:GLN:HG2	2:C:660:ASP:O	2.21	0.41
2:F:623:LEU:HA	2:F:699:ARG:HA	2.02	0.41
1:A:148:GLY:HA3	1:A:169:LEU:HD11	2.02	0.41
2:C:194:ASP:HB3	2:C:197:ASP:OD1	2.20	0.41
2:C:412:LEU:HD23	2:C:450:ILE:HD11	2.02	0.41
2:C:495:LEU:HD11	2:C:499:LEU:HD13	2.02	0.41
1:B:6:MET:HE2	1:B:34:GLN:CB	2.50	0.41
1:B:65:ASP:O	1:B:69:LYS:HB2	2.20	0.41
1:B:201:TYR:CE1	2:F:595:ALA:HB3	2.55	0.41
2:C:159:LEU:HD22	2:C:163:GLU:CB	2.45	0.41
2:C:405:LEU:CB	2:C:406:PRO:HD3	2.50	0.41
2:C:530:GLU:OE1	2:C:549:ARG:NH2	2.53	0.41
2:C:725:PRO:O	2:C:729:VAL:HG23	2.21	0.41
2:C:729:VAL:CG1	2:C:730:GLY:N	2.83	0.41
1:B:243:TYR:CZ	1:B:245:MET:HB2	2.55	0.41
2:F:485:HIS:O	2:F:592:HIS:N	2.49	0.41
1:B:64:SER:HB2	1:B:67:LYS:HB2	2.03	0.41
2:C:244:LEU:O	2:C:248:LEU:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:193:SER:O	1:B:195:VAL:HG23	2.21	0.41
2:C:390:GLU:HB2	2:C:694:MET:HE2	2.02	0.41
2:C:430:THR:HG21	2:C:432:GLU:OE1	2.21	0.41
2:F:553:TYR:O	2:F:554:LEU:CB	2.68	0.41
1:B:195:VAL:HB	1:B:208:ILE:HG22	2.03	0.40
1:B:275:LEU:HD22	1:B:284:PHE:CE1	2.56	0.40
2:C:44:GLY:HA3	2:C:245:SER:CB	2.52	0.40
2:C:411:VAL:O	2:C:415:VAL:HG23	2.21	0.40
2:C:623:LEU:HA	2:C:699:ARG:HA	2.03	0.40
2:C:653:VAL:HB	2:C:665:GLU:HB3	2.03	0.40
1:B:128:LEU:HD12	1:B:128:LEU:HA	1.96	0.40
2:C:117:VAL:HG21	2:C:133:LEU:HD13	2.02	0.40
2:C:442:HIS:O	2:C:446:ILE:HG12	2.21	0.40
2:C:93:ILE:HD11	2:C:357:ILE:CD1	2.52	0.40
2:C:176:ASN:OD1	2:C:193:VAL:HG22	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:217:GLN:OE1	2:C:217:GLN:OE1[3_655]	2.07	0.13

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	327/337 (97%)	315 (96%)	12 (4%)	0	100	100
1	B	323/337 (96%)	310 (96%)	13 (4%)	0	100	100
2	C	694/733 (95%)	665 (96%)	24 (4%)	5 (1%)	18	51
2	F	322/733 (44%)	310 (96%)	9 (3%)	3 (1%)	14	46

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1666/2140 (78%)	1600 (96%)	58 (4%)	8 (0%)	24 57

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	600	PRO
2	C	487	ARG
2	C	158	LYS
2	F	389	SER
2	C	387	HIS
2	C	539	TYR
2	F	594	ASP
2	C	72	ILE

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	282/299 (94%)	277 (98%)	5 (2%)	51 70
1	B	282/299 (94%)	272 (96%)	10 (4%)	32 58
2	C	583/621 (94%)	551 (94%)	32 (6%)	19 47
All	All	1147/1219 (94%)	1100 (96%)	47 (4%)	27 54

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	SER
1	A	128	LEU
1	A	151	LYS
1	A	327	ILE
1	A	328	LEU
1	B	68	MET
1	B	99	ASN
1	B	117	SER

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Mol	Chain	Res	Type
1	B	128	LEU
1	B	203	SER
1	B	208	ILE
1	B	222	ARG
1	B	236	VAL
1	B	276	LEU
1	B	304	LEU
2	C	31	ASP
2	C	36	THR
2	C	73	ASP
2	C	77	VAL
2	C	100	ASP
2	C	106	THR
2	C	147	ILE
2	C	172	TYR
2	C	235	GLN
2	C	242	VAL
2	C	248	LEU
2	C	267	VAL
2	C	277	SER
2	C	294	VAL
2	C	298	ASN
2	C	326	LEU
2	C	334	ARG
2	C	342	PHE
2	C	359	TYR
2	C	379	ILE
2	C	404	ASP
2	C	408	LEU
2	C	450	ILE
2	C	469	THR
2	C	492	VAL
2	C	493	GLU
2	C	499	LEU
2	C	536	TRP
2	C	594	ASP
2	C	645	ILE
2	C	665	GLU
2	C	666	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	GLN
1	A	157	GLN
1	A	269	ASN
1	B	127	HIS
1	B	129	HIS
1	B	152	ASN
1	B	157	GLN
1	B	162	ASN
2	C	81	HIS
2	C	82	ASN
2	C	208	ASN
2	C	255	HIS
2	C	337	GLN
2	C	348	ASN
2	C	458	GLN
2	C	476	GLN
2	C	504	GLN
2	C	727	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
3	SF4	B	401	1	0,12,12	-	-	-		
3	SF4	A	401	1	0,12,12	-	-	-		

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
3	SF4	B	401	1	-	-	0/6/5/5
3	SF4	A	401	1	-	-	0/6/5/5

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	401	SF4	1	0
3	A	401	SF4	3	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](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	329/337 (97%)	-0.58	2 (0%) 85 66	134, 205, 255, 285	0
1	B	325/337 (96%)	-0.62	0 100 100	124, 231, 302, 328	0
2	C	700/733 (95%)	-0.62	1 (0%) 92 88	87, 157, 218, 267	0
2	F	326/733 (44%)	-0.48	1 (0%) 90 77	275, 339, 391, 425	0
All	All	1680/2140 (78%)	-0.58	4 (0%) 91 83	87, 196, 361, 425	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	230	CYS	2.4
1	A	292	ILE	2.3
2	F	597	HIS	2.2
1	A	329	PHE	2.2

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SF4	A	401	8/8	0.97	0.03	154,168,236,240	0
3	SF4	B	401	8/8	0.99	0.03	164,225,278,279	0

6.5 Other polymers [i](#)

There are no such residues in this entry.