



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:36 PM UTC

PDB ID : 1MPY / pdb_00001mpy
Title : STRUCTURE OF CATECHOL 2,3-DIOXYGENASE (METAPYROCAT-
CHASE) FROM PSEUDOMONAS PUTIDA MT-2
Authors : Kita, A.; Kita, S.; Fujisawa, I.; Inaka, K.; Ishida, T.; Horiike, K.; Nozaki, M.;
Miki, K.
Deposited on : 1998-10-20
Resolution : 2.80 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

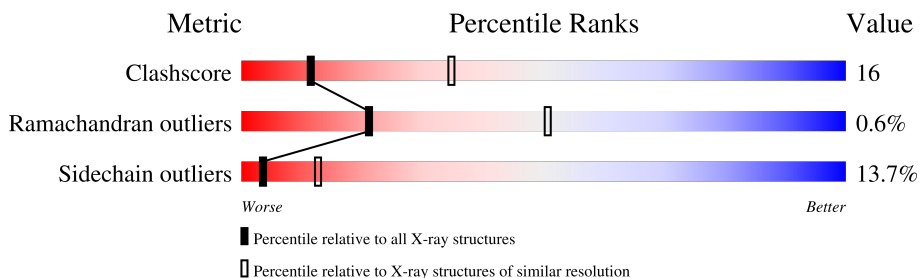
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	307	58% 33% 8% .
1	B	307	58% 36% 7%
1	C	307	57% 36% 7%
1	D	307	56% 37% 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10044 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.

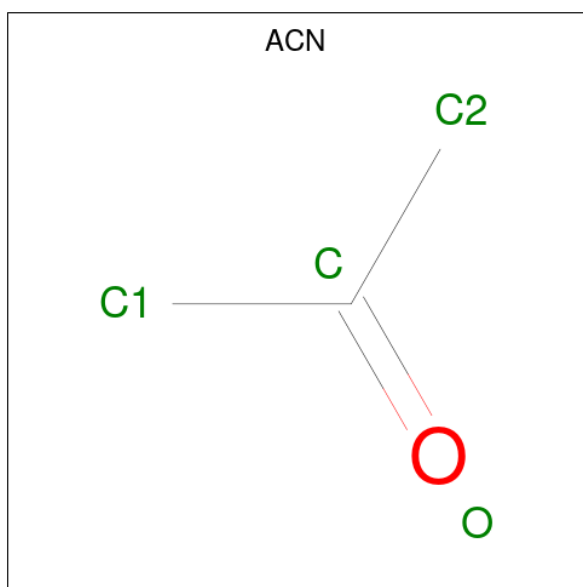
- Molecule 1 is a protein called CATECHOL 2,3-DIOXYGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	0	0
			2478	1574	430	460	14			
1	B	307	Total	C	N	O	S	0	0	0
			2478	1574	430	460	14			
1	C	307	Total	C	N	O	S	0	0	0
			2478	1574	430	460	14			
1	D	307	Total	C	N	O	S	0	0	0
			2478	1574	430	460	14			

- Molecule 2 is FE (II) ION (CCD ID: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Fe	0	0
			1	1		
2	B	1	Total	Fe	0	0
			1	1		
2	C	1	Total	Fe	0	0
			1	1		
2	D	1	Total	Fe	0	0
			1	1		

- Molecule 3 is ACETONE (CCD ID: ACN) (formula: C₃H₆O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	3	1		
3	B	1	Total	C	O	0	0
			4	3	1		
3	C	1	Total	C	O	0	0
			4	3	1		
3	D	1	Total	C	O	0	0
			4	3	1		

- Molecule 4 is water.

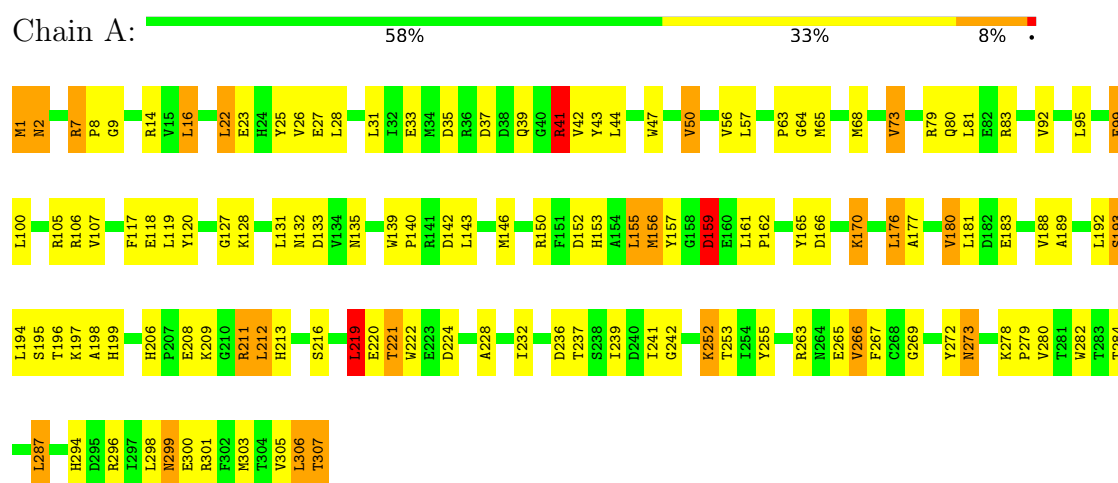
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total	O	0	0
			32	32		
4	B	24	Total	O	0	0
			24	24		
4	C	23	Total	O	0	0
			23	23		
4	D	33	Total	O	0	0
			33	33		

3 Residue-property plots

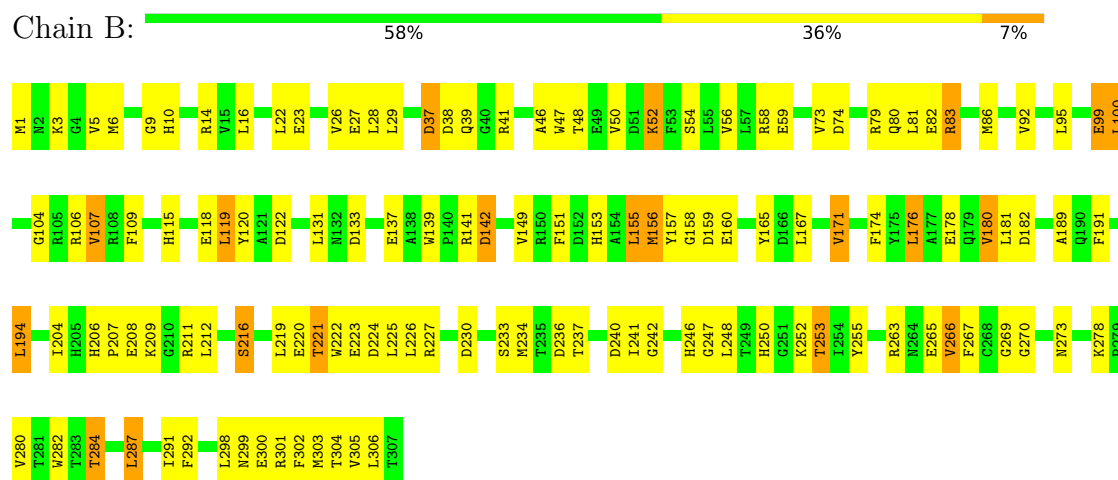
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

• Molecule 1: CATECHOL 2,3-DIOXYGENASE

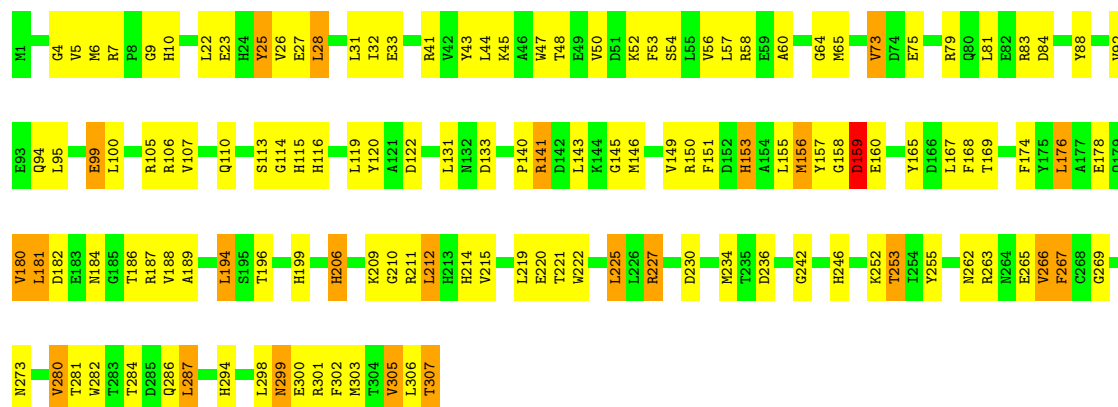


• Molecule 1: CATECHOL 2,3-DIOXYGENASE

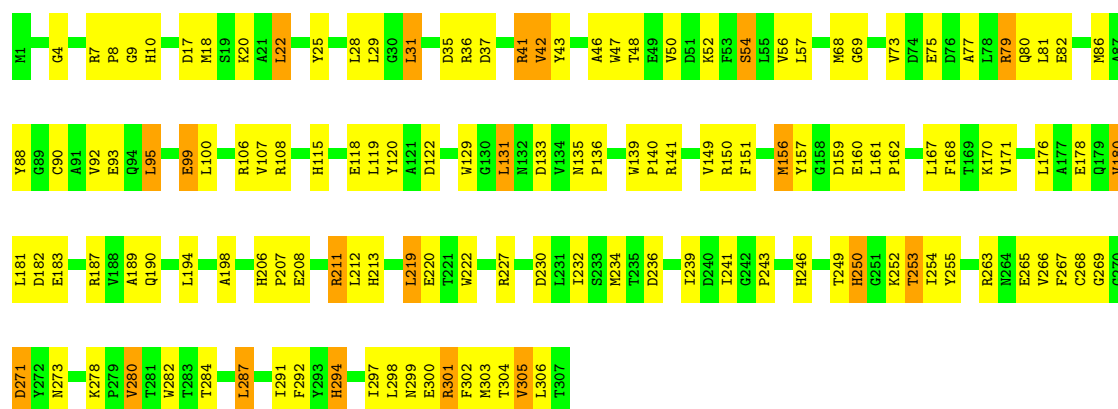


• Molecule 1: CATECHOL 2,3-DIOXYGENASE





• Molecule 1: CATECHOL 2,3-DIOXYGENASE



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	264.00Å 264.00Å 59.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.80	Depositor
% Data completeness (in resolution range)	84.0 (10.00-2.80)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.200 , 0.280	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	10044	wwPDB-VP
Average B, all atoms (Å ²)	14.0	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: FE2, ACN

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	0.58	0/2542	1.13	20/3440 (0.6%)
1	B	0.64	0/2542	1.16	23/3440 (0.7%)
1	C	0.64	1/2542 (0.0%)	1.16	26/3440 (0.8%)
1	D	0.62	1/2542 (0.0%)	1.15	20/3440 (0.6%)
All	All	0.62	2/10168 (0.0%)	1.15	89/13760 (0.6%)

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	C	0	2
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	141	ARG	C-N	6.37	1.42	1.33
1	D	271	ASP	C-N	-5.67	1.25	1.33

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	159	ASP	O-C-N	-12.19	107.81	122.19
1	D	159	ASP	N-CA-C	10.47	122.44	111.14
1	B	38	ASP	N-CA-C	-10.13	99.46	112.23
1	D	303	MET	CA-CB-CG	9.62	133.35	114.10
1	C	159	ASP	O-C-N	-9.52	108.30	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	47	TRP	N-CA-C	8.88	121.92	111.71
1	A	1	MET	CA-C-N	8.65	132.23	120.38
1	A	1	MET	C-N-CA	8.65	132.23	120.38
1	A	2	ASN	O-C-N	8.63	131.27	122.12
1	A	220	GLU	N-CA-C	8.41	120.37	111.03
1	B	142	ASP	N-CA-C	8.36	120.47	110.44
1	B	220	GLU	N-CA-C	7.91	121.44	111.24
1	C	47	TRP	N-CA-C	7.76	119.74	111.28
1	C	220	GLU	N-CA-C	7.67	119.55	111.03
1	D	303	MET	N-CA-CB	7.54	122.20	110.44
1	D	160	GLU	N-CA-C	7.53	119.80	109.11
1	B	47	TRP	N-CA-C	7.52	120.54	111.82
1	B	141	ARG	N-CA-C	7.37	122.10	113.18
1	D	141	ARG	N-CA-C	7.15	121.23	112.23
1	C	52	LYS	N-CA-C	7.11	120.86	111.75
1	B	216	SER	CB-CA-C	-7.03	98.21	109.80
1	C	151	PHE	N-CA-C	-7.03	99.26	109.59
1	A	253	THR	N-CA-C	7.01	119.83	109.24
1	C	305	VAL	N-CA-C	6.96	118.03	108.84
1	B	100	LEU	N-CA-C	-6.92	96.98	108.34
1	B	37	ASP	N-CA-C	6.92	120.62	110.17
1	D	168	PHE	N-CA-C	6.73	118.70	111.36
1	A	170	LYS	N-CA-C	6.63	118.19	110.97
1	C	73	VAL	N-CA-C	6.60	117.09	110.23
1	D	253	THR	N-CA-C	6.58	119.18	109.24
1	C	206	HIS	CA-C-N	6.58	126.52	119.28
1	C	206	HIS	C-N-CA	6.58	126.52	119.28
1	A	216	SER	N-CA-C	6.49	120.11	109.72
1	C	6	MET	N-CA-C	6.42	119.53	111.24
1	C	210	GLY	N-CA-C	6.40	123.54	115.42
1	A	155	LEU	CA-C-N	-6.31	113.48	122.94
1	A	155	LEU	C-N-CA	-6.31	113.48	122.94
1	C	25	TYR	N-CA-C	6.27	120.19	112.54
1	B	266	VAL	N-CA-C	-6.21	99.17	108.12
1	D	122	ASP	N-CA-C	6.19	118.48	109.07
1	B	160	GLU	CA-C-N	-6.10	106.91	121.80
1	B	160	GLU	C-N-CA	-6.10	106.91	121.80
1	A	41	ARG	N-CA-C	-6.07	101.50	110.48
1	D	303	MET	CB-CA-C	-6.02	97.33	109.55
1	A	269	GLY	N-CA-C	6.02	127.44	113.18
1	D	304	THR	N-CA-C	6.02	120.90	113.50
1	A	142	ASP	N-CA-C	6.01	120.42	112.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	294	HIS	N-CA-C	6.00	117.62	111.14
1	B	253	THR	N-CA-C	5.98	118.16	109.07
1	C	141	ARG	N-CA-C	5.89	120.20	112.89
1	A	99	GLU	N-CA-C	-5.77	104.18	111.11
1	D	139	TRP	CA-C-N	5.74	125.50	119.76
1	D	139	TRP	C-N-CA	5.74	125.50	119.76
1	B	160	GLU	O-C-N	5.73	131.57	122.81
1	B	141	ARG	CA-C-N	-5.66	113.16	123.34
1	B	141	ARG	C-N-CA	-5.66	113.16	123.34
1	A	278	LYS	N-CA-C	-5.63	102.95	109.93
1	D	88	TYR	N-CA-C	-5.62	105.69	112.54
1	D	305	VAL	N-CA-C	5.61	115.40	106.88
1	B	107	VAL	N-CA-C	-5.58	99.60	107.75
1	D	54	SER	N-CA-C	-5.57	108.28	114.62
1	B	247	GLY	N-CA-C	-5.56	105.76	112.49
1	C	176	LEU	CA-CB-CG	5.54	135.68	116.30
1	A	25	TYR	N-CA-C	5.51	118.21	111.82
1	C	267	PHE	N-CA-C	5.46	117.10	108.96
1	B	270	GLY	N-CA-C	-5.43	106.27	112.79
1	B	139	TRP	N-CA-C	5.42	113.00	108.13
1	B	151	PHE	N-CA-C	-5.41	101.36	109.15
1	A	152	ASP	N-CA-C	5.40	119.41	111.04
1	C	253	THR	N-CA-C	5.37	116.96	108.96
1	C	221	THR	N-CA-C	5.34	117.98	109.81
1	D	220	GLU	N-CA-C	5.34	119.73	112.04
1	A	221	THR	N-CA-C	5.33	117.97	109.81
1	D	211	ARG	N-CA-C	5.31	117.14	110.24
1	C	141	ARG	CA-C-N	-5.26	111.49	121.54
1	C	141	ARG	C-N-CA	-5.26	111.49	121.54
1	C	122	ASP	N-CA-C	5.25	116.79	108.96
1	C	242	GLY	CA-C-N	5.19	126.33	119.84
1	C	242	GLY	C-N-CA	5.19	126.33	119.84
1	D	46	ALA	N-CA-C	-5.08	102.26	110.14
1	D	170	LYS	N-CA-C	5.08	117.60	111.40
1	B	52	LYS	N-CA-C	5.07	117.19	111.11
1	B	221	THR	N-CA-C	5.06	117.56	109.81
1	B	46	ALA	N-CA-C	-5.06	102.30	110.14
1	C	153	HIS	CE1-NE2-CD2	-5.04	103.96	109.00
1	C	269	GLY	N-CA-C	5.04	125.11	113.18
1	C	266	VAL	CB-CA-C	5.02	117.89	110.77
1	C	215	VAL	N-CA-C	-5.01	101.26	108.48
1	A	219	LEU	N-CA-C	-5.00	101.24	109.40

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	159	ASP	Mainchain
1	C	159	ASP	Mainchain
1	C	160	GLU	Mainchain

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	2478	0	2385	73	0
1	B	2478	0	2385	81	0
1	C	2478	0	2385	90	0
1	D	2478	0	2385	83	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	4	0	6	0	0
3	B	4	0	6	0	0
3	C	4	0	6	0	0
3	D	4	0	6	0	0
4	A	32	0	0	2	0
4	B	24	0	0	0	0
4	C	23	0	0	0	0
4	D	33	0	0	1	0
All	All	10044	0	9564	313	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 16.

All (313) 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:C:180:VAL:HG13	1:C:189:ALA:HB3	1.57	0.86
1:C:64:GLY:HA2	1:C:307:THR:HG23	1.58	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:ILE:HG22	1:A:255:TYR:HB2	1.61	0.82
1:C:79:ARG:HG3	1:C:79:ARG:HH11	1.42	0.82
1:B:180:VAL:HG13	1:B:189:ALA:HB3	1.62	0.78
1:A:99:GLU:HB3	1:A:106:ARG:HH21	1.48	0.76
1:A:95:LEU:HD12	1:A:99:GLU:HG3	1.66	0.75
1:C:159:ASP:OD1	1:C:159:ASP:O	2.06	0.74
1:A:211:ARG:NH1	1:A:307:THR:HB	2.04	0.72
1:A:135:ASN:ND2	1:C:286:GLN:HG3	2.05	0.72
1:B:208:GLU:HG3	1:B:211:ARG:HD2	1.72	0.71
1:D:211:ARG:HE	1:D:211:ARG:HA	1.54	0.71
1:A:180:VAL:CG1	1:A:189:ALA:HB3	2.21	0.71
1:C:99:GLU:HG3	1:C:106:ARG:NH1	2.06	0.70
1:B:206:HIS:CD2	1:B:208:GLU:HG2	2.25	0.70
1:B:37:ASP:HB2	1:B:41:ARG:H	1.55	0.70
1:B:115:HIS:CD2	1:B:156:MET:HG2	2.26	0.69
1:D:115:HIS:CD2	1:D:156:MET:HG2	2.28	0.69
1:C:156:MET:HE3	1:C:212:LEU:HD23	1.75	0.69
1:D:180:VAL:CG1	1:D:189:ALA:HB3	2.21	0.69
1:A:140:PRO:HD2	1:A:143:LEU:HD11	1.75	0.68
1:A:219:LEU:HD22	1:A:266:VAL:HG11	1.77	0.66
1:B:99:GLU:HG3	1:B:106:ARG:NH1	2.11	0.66
1:D:25:TYR:HD1	1:D:29:LEU:HD12	1.60	0.66
1:C:149:VAL:HG12	1:C:150:ARG:HG2	1.77	0.66
1:D:77:ALA:O	1:D:81:LEU:HD23	1.96	0.66
1:C:25:TYR:O	1:C:31:LEU:HB2	1.95	0.65
1:D:95:LEU:HD21	1:D:108:ARG:HB2	1.78	0.65
1:C:178:GLU:HG2	1:C:280:VAL:HG13	1.77	0.65
1:B:5:VAL:HG12	1:B:194:LEU:HD21	1.79	0.64
1:D:187:ARG:HD3	1:D:190:GLN:HB3	1.80	0.64
1:A:9:GLY:HA3	1:A:120:TYR:OH	1.98	0.64
1:A:265:GLU:HG2	1:A:266:VAL:N	2.12	0.64
1:D:106:ARG:HG2	1:D:120:TYR:HB3	1.81	0.63
1:D:92:VAL:CG1	1:D:107:VAL:HG13	2.30	0.62
1:A:33:GLU:HA	1:A:44:LEU:HG	1.80	0.62
1:A:219:LEU:HD22	1:A:266:VAL:CG1	2.29	0.62
1:D:10:HIS:HA	1:D:54:SER:O	2.00	0.62
1:B:302:PHE:CE2	1:B:303:MET:HG2	2.35	0.62
1:D:48:THR:HG21	1:D:149:VAL:HG13	1.81	0.62
1:B:99:GLU:HG3	1:B:106:ARG:HH12	1.65	0.62
1:A:211:ARG:HB3	1:A:211:ARG:HH11	1.65	0.61
1:C:28:LEU:O	1:C:145:GLY:HA3	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:ARG:HA	1:B:86:MET:HE3	1.83	0.61
1:C:227:ARG:HG2	1:C:227:ARG:HH11	1.66	0.60
1:A:99:GLU:HB3	1:A:106:ARG:NH2	2.15	0.60
1:C:156:MET:CE	1:C:212:LEU:HD23	2.32	0.59
1:B:227:ARG:HG2	1:B:227:ARG:HH11	1.67	0.59
1:B:284:THR:HA	1:B:287:LEU:HD23	1.84	0.59
1:A:211:ARG:HH12	1:A:306:LEU:HB3	1.68	0.59
1:C:252:LYS:O	1:C:267:PHE:HB2	2.03	0.59
1:D:301:ARG:H	1:D:301:ARG:HD3	1.68	0.59
1:B:23:GLU:O	1:B:27:GLU:HB2	2.03	0.59
1:A:252:LYS:O	1:A:267:PHE:HB2	2.02	0.58
1:D:95:LEU:HD23	1:D:106:ARG:HB2	1.85	0.58
1:B:3:LYS:HB3	1:B:81:LEU:HD21	1.84	0.58
1:D:25:TYR:CD1	1:D:29:LEU:HD12	2.38	0.58
1:B:37:ASP:HB3	1:B:39:GLN:H	1.68	0.58
1:B:159:ASP:HB3	1:B:209:LYS:HZ3	1.68	0.58
1:B:10:HIS:HA	1:B:54:SER:O	2.04	0.58
1:C:79:ARG:O	1:C:83:ARG:HG2	2.04	0.57
1:C:113:SER:HB2	1:C:158:GLY:HA3	1.86	0.57
1:D:82:GLU:O	1:D:86:MET:HG3	2.04	0.57
1:A:1:MET:HE3	1:A:195:SER:HB2	1.86	0.57
1:B:9:GLY:HA3	1:B:120:TYR:OH	2.05	0.56
1:C:65:MET:SD	1:C:156:MET:HE2	2.45	0.56
1:B:159:ASP:HB3	1:B:209:LYS:NZ	2.20	0.56
1:A:180:VAL:HA	1:A:282:TRP:O	2.06	0.56
1:C:23:GLU:O	1:C:27:GLU:HB2	2.06	0.56
1:D:93:GLU:HB3	1:D:108:ARG:HB3	1.88	0.56
1:A:211:ARG:NH1	1:A:306:LEU:HB3	2.20	0.55
1:B:100:LEU:HB2	1:B:106:ARG:NH1	2.21	0.55
1:C:176:LEU:HD21	1:C:187:ARG:HE	1.70	0.55
1:C:246:HIS:NE2	1:C:253:THR:HG21	2.21	0.55
1:C:253:THR:OG1	1:C:265:GLU:HG3	2.06	0.55
1:D:161:LEU:HB2	1:D:162:PRO:HD3	1.88	0.55
1:D:287:LEU:HD12	1:D:298:LEU:HD11	1.89	0.55
1:C:43:TYR:HE2	1:C:100:LEU:HD22	1.72	0.55
1:A:180:VAL:HG12	1:A:189:ALA:HB3	1.89	0.55
1:D:25:TYR:O	1:D:31:LEU:HB2	2.07	0.55
1:B:301:ARG:HA	1:B:304:THR:OG1	2.06	0.55
1:D:263:ARG:NH2	1:D:305:VAL:O	2.39	0.54
1:D:301:ARG:HH11	1:D:301:ARG:HB2	1.73	0.54
1:B:219:LEU:HD23	1:B:225:LEU:HD12	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:92:VAL:CG1	1:B:107:VAL:HG13	2.38	0.54
1:B:109:PHE:HE1	1:B:119:LEU:HD13	1.73	0.54
1:D:99:GLU:HG3	1:D:106:ARG:NH2	2.23	0.54
1:D:182:ASP:HA	1:D:284:THR:OG1	2.08	0.54
1:C:99:GLU:HG3	1:C:106:ARG:HH12	1.73	0.54
1:A:211:ARG:NH1	1:A:211:ARG:HB3	2.23	0.54
1:D:255:TYR:CE1	1:D:265:GLU:HB2	2.43	0.54
1:C:182:ASP:HA	1:C:284:THR:OG1	2.07	0.54
1:C:206:HIS:CD2	1:C:211:ARG:HG3	2.43	0.54
1:C:153:HIS:CD2	1:C:199:HIS:NE2	2.76	0.54
1:A:206:HIS:CD2	1:A:208:GLU:HB2	2.43	0.53
1:D:180:VAL:HG12	1:D:189:ALA:HB3	1.90	0.53
1:C:4:GLY:O	1:C:73:VAL:HG23	2.09	0.52
1:D:219:LEU:HD22	1:D:266:VAL:HG22	1.90	0.52
1:C:79:ARG:HG3	1:C:79:ARG:NH1	2.12	0.52
1:A:23:GLU:O	1:A:27:GLU:HB2	2.10	0.52
1:A:153:HIS:CD2	1:A:199:HIS:NE2	2.77	0.52
1:A:37:ASP:OD2	1:A:41:ARG:HD3	2.09	0.52
1:B:137:GLU:HB3	1:C:234:MET:HE2	1.92	0.52
1:B:300:GLU:HG3	1:B:301:ARG:HG3	1.90	0.52
1:D:92:VAL:HG13	1:D:107:VAL:HG13	1.91	0.52
1:A:166:ASP:O	1:A:170:LYS:HB2	2.11	0.51
1:C:252:LYS:HD3	1:C:252:LYS:N	2.25	0.51
1:C:5:VAL:HG12	1:C:194:LEU:HD21	1.92	0.51
1:A:299:ASN:HD21	1:A:301:ARG:HB2	1.75	0.51
1:B:180:VAL:CG1	1:B:189:ALA:HB3	2.37	0.51
1:D:178:GLU:HG2	1:D:280:VAL:HG13	1.92	0.51
1:A:284:THR:O	1:A:287:LEU:HB2	2.10	0.51
1:B:22:LEU:O	1:B:26:VAL:HG22	2.10	0.51
1:C:41:ARG:NH2	1:C:100:LEU:HD23	2.26	0.51
1:B:92:VAL:HG13	1:B:107:VAL:HG13	1.93	0.51
1:D:106:ARG:HD2	1:D:118:GLU:OE2	2.11	0.51
1:D:252:LYS:O	1:D:267:PHE:HB2	2.10	0.51
1:D:299:ASN:HD21	1:D:301:ARG:NH1	2.08	0.51
1:A:222:TRP:NE1	1:B:226:LEU:HB2	2.26	0.51
1:A:263:ARG:HH22	1:A:305:VAL:HB	1.75	0.51
1:A:232:ILE:HG23	1:A:237:THR:HB	1.92	0.50
1:C:230:ASP:O	1:C:234:MET:HG3	2.11	0.50
1:A:8:PRO:HB3	1:A:68:MET:HE3	1.93	0.50
1:A:79:ARG:HG3	1:A:105:ARG:NH2	2.27	0.50
1:A:146:MET:HE2	1:A:228:ALA:HB1	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:HIS:O	1:A:296:ARG:NH1	2.44	0.50
1:B:14:ARG:HG2	1:B:58:ARG:NH1	2.27	0.50
1:B:246:HIS:CE1	1:B:253:THR:HG21	2.46	0.50
1:B:300:GLU:HG3	1:B:301:ARG:H	1.76	0.50
1:B:191:PHE:CE1	1:B:248:LEU:HD13	2.46	0.50
1:B:1:MET:N	1:B:194:LEU:O	2.44	0.50
1:C:146:MET:HA	1:C:227:ARG:NH1	2.27	0.50
1:D:92:VAL:HG11	1:D:107:VAL:HG13	1.92	0.50
1:C:294:HIS:HE1	1:D:230:ASP:OD1	1.93	0.50
1:D:255:TYR:CD1	1:D:265:GLU:HB2	2.47	0.50
1:A:92:VAL:HG13	1:A:107:VAL:HG13	1.93	0.50
1:B:287:LEU:HD12	1:B:298:LEU:HD11	1.94	0.50
1:D:99:GLU:HG3	1:D:106:ARG:HH21	1.75	0.50
1:B:299:ASN:O	1:B:302:PHE:HB3	2.11	0.49
1:C:41:ARG:HH21	1:C:100:LEU:HD23	1.77	0.49
1:C:75:GLU:O	1:C:79:ARG:HG2	2.11	0.49
1:D:291:ILE:HG21	1:D:302:PHE:CE2	2.48	0.49
1:B:241:ILE:HD13	1:B:292:PHE:CE1	2.48	0.49
1:C:32:ILE:N	1:C:32:ILE:HD12	2.27	0.49
1:D:211:ARG:HA	1:D:211:ARG:NE	2.24	0.49
1:B:155:LEU:HD21	1:B:204:ILE:HG13	1.95	0.49
1:B:165:TYR:HE1	1:B:176:LEU:HG	1.77	0.49
1:C:48:THR:HG21	1:C:149:VAL:HG13	1.95	0.49
1:D:99:GLU:CG	1:D:106:ARG:NH2	2.75	0.49
1:D:131:LEU:HD13	1:D:140:PRO:HG3	1.94	0.49
1:C:301:ARG:HG2	1:C:305:VAL:HG21	1.95	0.49
1:D:278:LYS:NZ	1:D:278:LYS:HB3	2.28	0.49
1:C:255:TYR:CE1	1:C:265:GLU:HB2	2.48	0.48
1:A:22:LEU:HD21	1:A:42:VAL:HG11	1.94	0.48
1:B:157:TYR:HB2	1:B:211:ARG:HB2	1.94	0.48
1:B:246:HIS:NE2	1:B:253:THR:HG21	2.28	0.48
1:B:252:LYS:O	1:B:267:PHE:HB2	2.14	0.48
1:D:18:MET:SD	1:D:42:VAL:HG12	2.53	0.48
1:C:114:GLY:HA3	1:C:209:LYS:HB3	1.95	0.48
1:A:161:LEU:HB2	1:A:162:PRO:HD3	1.93	0.48
1:A:117:PHE:CE1	1:A:156:MET:HE3	2.48	0.48
1:D:8:PRO:HB3	1:D:68:MET:HE2	1.96	0.48
1:D:246:HIS:O	1:D:250:HIS:HA	2.13	0.48
1:B:298:LEU:HD22	1:B:302:PHE:CD2	2.49	0.48
1:C:299:ASN:O	1:C:302:PHE:HB3	2.12	0.48
1:C:262:ASN:OD1	1:C:307:THR:HG21	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:ASP:OD2	1:B:41:ARG:HD3	2.14	0.48
1:C:9:GLY:HA3	1:C:120:TYR:OH	2.13	0.48
1:C:33:GLU:HA	1:C:44:LEU:HG	1.96	0.48
1:C:181:LEU:HD12	1:C:187:ARG:HA	1.95	0.48
1:D:284:THR:O	1:D:287:LEU:HB2	2.14	0.47
1:C:10:HIS:HA	1:C:54:SER:O	2.14	0.47
1:D:301:ARG:HA	1:D:305:VAL:HG23	1.95	0.47
1:B:37:ASP:HB2	1:B:41:ARG:N	2.27	0.47
1:D:246:HIS:CE1	1:D:253:THR:HG21	2.49	0.47
1:A:193:SER:HB3	1:A:197:LYS:O	2.14	0.47
1:B:174:PHE:CD2	1:B:194:LEU:HD13	2.49	0.47
1:D:75:GLU:HG3	1:D:79:ARG:NH1	2.30	0.47
1:B:178:GLU:HB2	1:B:191:PHE:HB2	1.96	0.47
1:C:157:TYR:HE2	1:C:303:MET:SD	2.37	0.47
1:C:263:ARG:NH2	1:C:305:VAL:O	2.47	0.47
1:A:65:MET:HB2	1:A:212:LEU:HG	1.96	0.47
1:D:22:LEU:HD12	1:D:22:LEU:HA	1.74	0.47
1:D:180:VAL:HA	1:D:282:TRP:O	2.14	0.47
1:D:219:LEU:HD12	1:D:219:LEU:HA	1.70	0.47
1:B:263:ARG:NH2	1:B:305:VAL:O	2.48	0.46
1:C:64:GLY:HA2	1:C:307:THR:CG2	2.37	0.46
1:C:92:VAL:HG13	1:C:107:VAL:HG13	1.97	0.46
1:C:168:PHE:O	1:C:174:PHE:HB2	2.15	0.46
1:D:227:ARG:HD3	4:D:5104:HOH:O	2.13	0.46
1:B:255:TYR:CE1	1:B:265:GLU:HB2	2.51	0.46
1:A:296:ARG:HG3	1:A:296:ARG:HH11	1.80	0.46
1:B:5:VAL:CG1	1:B:194:LEU:HD21	2.45	0.46
1:C:28:LEU:HD22	1:C:146:MET:HE3	1.98	0.46
1:D:35:ASP:O	1:D:36:ARG:HG3	2.15	0.46
1:B:302:PHE:CD2	1:B:303:MET:HG2	2.51	0.46
1:D:18:MET:SD	1:D:42:VAL:CG1	3.04	0.46
1:A:16:LEU:HD11	1:A:63:PRO:HD3	1.98	0.46
1:A:50:VAL:HG22	1:A:127:GLY:HA2	1.98	0.46
1:B:79:ARG:HG2	1:B:79:ARG:HH11	1.81	0.46
1:D:18:MET:O	1:D:22:LEU:HB2	2.15	0.46
1:A:106:ARG:HD2	1:A:118:GLU:OE1	2.16	0.46
1:C:79:ARG:HD3	1:C:105:ARG:NH2	2.31	0.46
1:C:287:LEU:HD13	1:C:298:LEU:HD11	1.98	0.46
1:A:35:ASP:HB3	1:A:43:TYR:CD1	2.51	0.45
1:B:104:GLY:HA3	1:B:122:ASP:HB2	1.97	0.45
1:C:212:LEU:HB3	1:C:307:THR:HG22	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:291:ILE:HD12	1:D:302:PHE:CZ	2.51	0.45
1:A:299:ASN:ND2	1:A:299:ASN:C	2.74	0.45
1:D:135:ASN:N	1:D:136:PRO:CD	2.78	0.45
1:A:221:THR:O	1:A:224:ASP:HB2	2.16	0.45
1:C:159:ASP:OD1	1:C:159:ASP:C	2.59	0.45
1:D:232:ILE:HG21	1:D:239:ILE:HB	1.98	0.45
1:B:82:GLU:O	1:B:86:MET:HG3	2.17	0.45
1:B:153:HIS:NE2	1:B:216:SER:OG	2.50	0.45
1:D:252:LYS:HD2	1:D:268:CYS:SG	2.57	0.45
1:A:159:ASP:HB3	1:A:209:LYS:HE2	1.99	0.45
1:B:107:VAL:O	1:B:118:GLU:HA	2.17	0.45
1:B:300:GLU:HG3	1:B:301:ARG:N	2.31	0.45
1:B:180:VAL:HA	1:B:282:TRP:O	2.16	0.45
1:D:9:GLY:O	1:D:54:SER:HA	2.17	0.45
1:A:177:ALA:O	1:A:279:PRO:HA	2.17	0.45
1:A:294:HIS:HE1	1:B:230:ASP:OD1	2.00	0.45
1:D:17:ASP:OD2	1:D:20:LYS:HG3	2.16	0.45
1:D:99:GLU:CG	1:D:106:ARG:HH21	2.29	0.45
1:C:43:TYR:C	1:C:44:LEU:HD12	2.42	0.44
1:C:255:TYR:CD1	1:C:265:GLU:HB2	2.51	0.44
1:D:4:GLY:O	1:D:73:VAL:HG23	2.17	0.44
1:D:181:LEU:HD23	1:D:187:ARG:HA	1.99	0.44
1:A:128:LYS:HE2	4:A:5012:HOH:O	2.17	0.44
1:C:84:ASP:O	1:C:88:TYR:HB2	2.18	0.44
1:D:206:HIS:ND1	1:D:207:PRO:HD2	2.33	0.44
1:B:48:THR:HG21	1:B:149:VAL:HG13	1.98	0.44
1:C:92:VAL:CG1	1:C:107:VAL:HG13	2.48	0.44
1:A:7:ARG:HD2	1:A:7:ARG:N	2.32	0.44
1:A:294:HIS:HA	1:B:233:SER:OG	2.17	0.44
1:C:65:MET:SD	1:C:156:MET:CE	3.05	0.44
1:A:135:ASN:OD1	1:C:282:TRP:HA	2.17	0.44
1:A:252:LYS:HE2	1:B:226:LEU:HD21	2.00	0.44
1:A:132:ASN:OD1	1:A:133:ASP:N	2.51	0.44
1:A:252:LYS:N	1:A:252:LYS:CD	2.80	0.44
1:A:272:TYR:CZ	1:C:150:ARG:HD3	2.53	0.44
1:B:158:GLY:O	1:B:206:HIS:N	2.44	0.44
1:C:165:TYR:CE1	1:C:169:THR:HG21	2.54	0.43
1:C:196:THR:HA	1:C:273:ASN:O	2.18	0.43
1:C:222:TRP:CD1	1:D:222:TRP:CD1	3.06	0.43
1:A:198:ALA:O	1:A:199:HIS:HB3	2.17	0.43
1:D:250:HIS:HB2	1:D:271:ASP:OD2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:254:ILE:HB	1:D:266:VAL:HG12	1.99	0.43
1:C:176:LEU:HD21	1:C:187:ARG:NE	2.32	0.43
1:C:255:TYR:HE1	1:C:265:GLU:CD	2.26	0.43
1:D:206:HIS:CD2	1:D:208:GLU:HG2	2.53	0.43
1:B:167:LEU:HD23	1:B:171:VAL:HG13	1.99	0.43
1:D:68:MET:HE3	1:D:151:PHE:HZ	1.83	0.43
1:C:95:LEU:N	1:C:95:LEU:HD22	2.33	0.43
1:D:292:PHE:HD2	1:D:297:ILE:O	2.02	0.43
1:D:47:TRP:HB2	1:D:129:TRP:HB3	2.00	0.43
1:B:37:ASP:HB3	1:B:39:GLN:N	2.33	0.43
1:D:198:ALA:HB1	1:D:249:THR:HG22	2.01	0.43
1:A:139:TRP:CE2	1:D:234:MET:HE1	2.54	0.43
1:C:214:HIS:HB3	1:C:263:ARG:HG2	2.00	0.43
1:D:206:HIS:CD2	1:D:211:ARG:HG3	2.54	0.43
1:C:140:PRO:HG2	1:C:143:LEU:HD11	2.00	0.43
1:A:165:TYR:HE1	1:A:176:LEU:HG	1.84	0.42
1:A:14:ARG:NH2	1:A:64:GLY:HA3	2.34	0.42
1:D:69:GLY:HA2	1:D:118:GLU:O	2.19	0.42
1:C:214:HIS:CB	1:C:263:ARG:HG2	2.49	0.42
1:C:230:ASP:OD1	1:D:294:HIS:HE1	2.03	0.42
1:C:214:HIS:C	1:C:214:HIS:CD2	2.97	0.42
1:D:9:GLY:HA3	1:D:120:TYR:OH	2.20	0.42
1:A:296:ARG:NH2	1:B:237:THR:O	2.52	0.42
1:B:6:MET:HE3	1:B:6:MET:HB2	1.86	0.42
1:B:216:SER:OG	1:B:265:GLU:OE1	2.30	0.42
1:D:22:LEU:HD11	1:D:42:VAL:HG11	2.02	0.42
1:C:159:ASP:HB3	1:C:209:LYS:HZ3	1.85	0.42
1:D:37:ASP:OD2	1:D:41:ARG:HD3	2.20	0.42
1:D:157:TYR:CE1	1:D:213:HIS:HD2	2.38	0.42
1:A:157:TYR:CE1	1:A:213:HIS:CD2	3.08	0.42
1:C:110:GLN:HB2	1:C:116:HIS:CE1	2.55	0.41
1:A:106:ARG:HG2	1:A:120:TYR:HB3	2.02	0.41
1:C:58:ARG:NH1	1:C:60:ALA:HB2	2.34	0.41
1:B:194:LEU:HD12	1:B:194:LEU:HA	1.86	0.41
1:A:7:ARG:NH1	4:A:5001:HOH:O	2.54	0.41
1:A:196:THR:HA	1:A:273:ASN:O	2.21	0.41
1:A:222:TRP:CD1	1:B:222:TRP:CD1	3.09	0.41
1:A:157:TYR:OH	1:A:303:MET:HG3	2.20	0.41
1:C:79:ARG:NH1	1:C:79:ARG:CG	2.82	0.41
1:B:37:ASP:OD2	1:B:41:ARG:HB2	2.20	0.41
1:B:182:ASP:HA	1:B:284:THR:HB	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:206:HIS:CD2	1:B:208:GLU:CG	3.02	0.41
1:B:206:HIS:HA	1:B:207:PRO:HD3	1.93	0.41
1:C:65:MET:HB2	1:C:212:LEU:HG	2.02	0.41
1:C:100:LEU:HB2	1:C:106:ARG:NH1	2.36	0.41
1:C:115:HIS:CD2	1:C:156:MET:HG3	2.55	0.41
1:C:227:ARG:HH11	1:C:227:ARG:CG	2.34	0.41
1:D:43:TYR:CD1	1:D:43:TYR:N	2.87	0.41
1:D:100:LEU:HD23	1:D:100:LEU:HA	1.89	0.41
1:A:100:LEU:HD23	1:A:100:LEU:HA	1.91	0.40
1:B:291:ILE:HG13	1:B:298:LEU:HD21	2.03	0.40
1:A:43:TYR:C	1:A:44:LEU:HD12	2.45	0.40
1:B:221:THR:O	1:B:224:ASP:HB2	2.20	0.40
1:B:240:ASP:HB3	1:B:255:TYR:HB3	2.02	0.40
1:B:246:HIS:O	1:B:250:HIS:HA	2.20	0.40
1:C:181:LEU:HD22	1:C:281:THR:CG2	2.51	0.40
1:C:181:LEU:HD22	1:C:281:THR:HG23	2.04	0.40
1:C:222:TRP:O	1:C:225:LEU:HB2	2.21	0.40
1:B:14:ARG:HG2	1:B:58:ARG:HH12	1.86	0.40
1:C:45:LYS:HB3	1:C:53:PHE:HA	2.03	0.40
1:C:182:ASP:HB3	1:C:188:VAL:HG21	2.03	0.40
1:A:239:ILE:HD11	1:A:242:GLY:HA2	2.04	0.40
1:C:186:THR:O	1:C:188:VAL:HG23	2.21	0.40
1:D:47:TRP:CZ3	1:D:140:PRO:HD3	2.57	0.40
1:B:73:VAL:HG23	1:B:74:ASP:H	1.86	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	305/307 (99%)	281 (92%)	23 (8%)	1 (0%)	36 66

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	305/307 (99%)	290 (95%)	13 (4%)	2 (1%)	18	47
1	C	305/307 (99%)	286 (94%)	18 (6%)	1 (0%)	36	66
1	D	305/307 (99%)	288 (94%)	14 (5%)	3 (1%)	12	38
All	All	1220/1228 (99%)	1145 (94%)	68 (6%)	7 (1%)	21	51

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	300	GLU
1	B	242	GLY
1	B	269	GLY
1	D	250	HIS
1	A	73	VAL
1	D	243	PRO
1	D	269	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 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	262/262 (100%)	219 (84%)	43 (16%)	2	8
1	B	262/262 (100%)	229 (87%)	33 (13%)	4	15
1	C	262/262 (100%)	230 (88%)	32 (12%)	5	16
1	D	262/262 (100%)	226 (86%)	36 (14%)	3	13
All	All	1048/1048 (100%)	904 (86%)	144 (14%)	3	13

All (144) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	7	ARG
1	A	16	LEU
1	A	22	LEU

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Mol	Chain	Res	Type
1	A	26	VAL
1	A	28	LEU
1	A	31	LEU
1	A	39	GLN
1	A	41	ARG
1	A	50	VAL
1	A	56	VAL
1	A	57	LEU
1	A	73	VAL
1	A	80	GLN
1	A	81	LEU
1	A	83	ARG
1	A	119	LEU
1	A	131	LEU
1	A	150	ARG
1	A	155	LEU
1	A	156	MET
1	A	176	LEU
1	A	180	VAL
1	A	181	LEU
1	A	183	GLU
1	A	188	VAL
1	A	192	LEU
1	A	193	SER
1	A	194	LEU
1	A	211	ARG
1	A	212	LEU
1	A	219	LEU
1	A	236	ASP
1	A	252	LYS
1	A	266	VAL
1	A	273	ASN
1	A	280	VAL
1	A	287	LEU
1	A	298	LEU
1	A	299	ASN
1	A	300	GLU
1	A	306	LEU
1	A	307	THR
1	B	16	LEU
1	B	28	LEU
1	B	29	LEU

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Mol	Chain	Res	Type
1	B	50	VAL
1	B	52	LYS
1	B	56	VAL
1	B	59	GLU
1	B	80	GLN
1	B	83	ARG
1	B	95	LEU
1	B	99	GLU
1	B	119	LEU
1	B	131	LEU
1	B	133	ASP
1	B	142	ASP
1	B	155	LEU
1	B	156	MET
1	B	171	VAL
1	B	176	LEU
1	B	180	VAL
1	B	181	LEU
1	B	194	LEU
1	B	212	LEU
1	B	223	GLU
1	B	234	MET
1	B	236	ASP
1	B	266	VAL
1	B	273	ASN
1	B	278	LYS
1	B	280	VAL
1	B	284	THR
1	B	287	LEU
1	B	306	LEU
1	C	7	ARG
1	C	22	LEU
1	C	26	VAL
1	C	28	LEU
1	C	50	VAL
1	C	56	VAL
1	C	57	LEU
1	C	81	LEU
1	C	94	GLN
1	C	99	GLU
1	C	119	LEU
1	C	131	LEU

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Mol	Chain	Res	Type
1	C	133	ASP
1	C	141	ARG
1	C	155	LEU
1	C	156	MET
1	C	167	LEU
1	C	180	VAL
1	C	181	LEU
1	C	184	ASN
1	C	194	LEU
1	C	212	LEU
1	C	219	LEU
1	C	225	LEU
1	C	227	ARG
1	C	236	ASP
1	C	266	VAL
1	C	280	VAL
1	C	287	LEU
1	C	299	ASN
1	C	306	LEU
1	C	307	THR
1	D	7	ARG
1	D	22	LEU
1	D	28	LEU
1	D	31	LEU
1	D	41	ARG
1	D	42	VAL
1	D	50	VAL
1	D	52	LYS
1	D	56	VAL
1	D	57	LEU
1	D	79	ARG
1	D	80	GLN
1	D	90	CYS
1	D	95	LEU
1	D	99	GLU
1	D	119	LEU
1	D	131	LEU
1	D	133	ASP
1	D	150	ARG
1	D	156	MET
1	D	167	LEU
1	D	171	VAL

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Mol	Chain	Res	Type
1	D	176	LEU
1	D	180	VAL
1	D	183	GLU
1	D	194	LEU
1	D	212	LEU
1	D	219	LEU
1	D	236	ASP
1	D	241	ILE
1	D	273	ASN
1	D	280	VAL
1	D	287	LEU
1	D	300	GLU
1	D	301	ARG
1	D	306	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	39	GLN
1	A	206	HIS
1	A	218	HIS
1	A	277	HIS
1	A	294	HIS
1	A	299	ASN
1	B	110	GLN
1	B	184	ASN
1	B	199	HIS
1	B	206	HIS
1	C	2	ASN
1	C	12	GLN
1	C	94	GLN
1	C	206	HIS
1	C	277	HIS
1	C	294	HIS
1	D	116	HIS
1	D	218	HIS
1	D	294	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](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	$\# Z > 2$
3	ACN	B	309	2	3,3,3	0.50	0	3,3,3	0.66	0
3	ACN	D	309	2	3,3,3	1.22	0	3,3,3	0.93	0
3	ACN	C	309	2	3,3,3	0.47	0	3,3,3	0.52	0
3	ACN	A	309	2	3,3,3	0.87	0	3,3,3	0.70	0

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.

No monomer is involved in short contacts.

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.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.