



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:47 PM GMT

PDB ID : 1YNI
Title : Crystal Structure of N-Succinylarginine Dihydrolase, AstB, bound to Substrate and Product, an Enzyme from the Arginine Catabolic Pathway of Escherichia coli
Authors : Tocilj, A.; Schrag, J.D.; Li, Y.; Schneider, B.L.; Reitzer, L.; Matte, A.; Cygler, M.; Montreal-Kingston Bacterial Structural Genomics Initiative (BSGI)
Deposited on : 2005-01-24
Resolution : 2.20 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) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

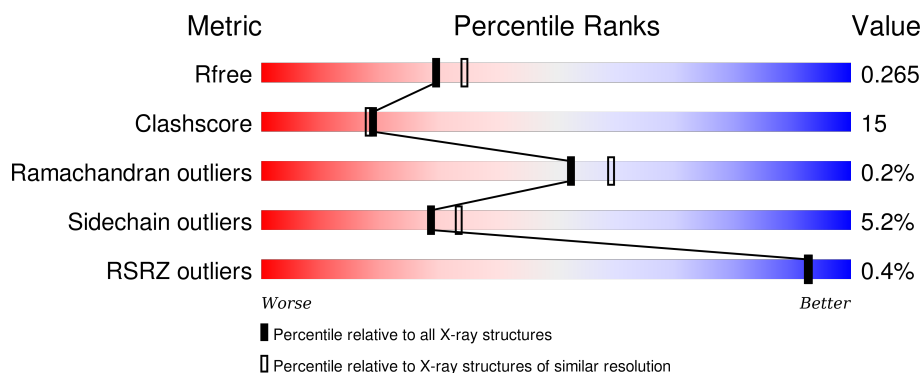
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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.

Mol	Chain	Length	Quality of chain
1	A	458	<div> <div>68%</div> <div>27%</div> <div>• •</div> </div>
1	B	458	<div> <div>71%</div> <div>22%</div> <div>• •</div> </div>
1	C	458	<div> <div>%</div> <div>66%</div> <div>27%</div> <div>• •</div> </div>
1	D	458	<div> <div>66%</div> <div>28%</div> <div>• •</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14363 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 Succinylarginine Dihydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	0	0
			3429	2148	626	644	11			
1	B	440	Total	C	N	O	S	0	0	0
			3429	2148	626	644	11			
1	C	440	Total	C	N	O	S	0	0	0
			3429	2148	626	644	11			
1	D	440	Total	C	N	O	S	0	0	0
			3429	2148	626	644	11			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	EXPRESSION TAG	UNP P76216
A	-9	GLY	-	EXPRESSION TAG	UNP P76216
A	-8	SER	-	EXPRESSION TAG	UNP P76216
A	-7	SER	-	EXPRESSION TAG	UNP P76216
A	-6	HIS	-	EXPRESSION TAG	UNP P76216
A	-5	HIS	-	EXPRESSION TAG	UNP P76216
A	-4	HIS	-	EXPRESSION TAG	UNP P76216
A	-3	HIS	-	EXPRESSION TAG	UNP P76216
A	-2	HIS	-	EXPRESSION TAG	UNP P76216
A	-1	HIS	-	EXPRESSION TAG	UNP P76216
A	0	GLY	-	EXPRESSION TAG	UNP P76216
A	1	SER	-	EXPRESSION TAG	UNP P76216
A	365	SER	CYS	ENGINEERED	UNP P76216
B	-10	MET	-	EXPRESSION TAG	UNP P76216
B	-9	GLY	-	EXPRESSION TAG	UNP P76216
B	-8	SER	-	EXPRESSION TAG	UNP P76216
B	-7	SER	-	EXPRESSION TAG	UNP P76216
B	-6	HIS	-	EXPRESSION TAG	UNP P76216
B	-5	HIS	-	EXPRESSION TAG	UNP P76216
B	-4	HIS	-	EXPRESSION TAG	UNP P76216
B	-3	HIS	-	EXPRESSION TAG	UNP P76216

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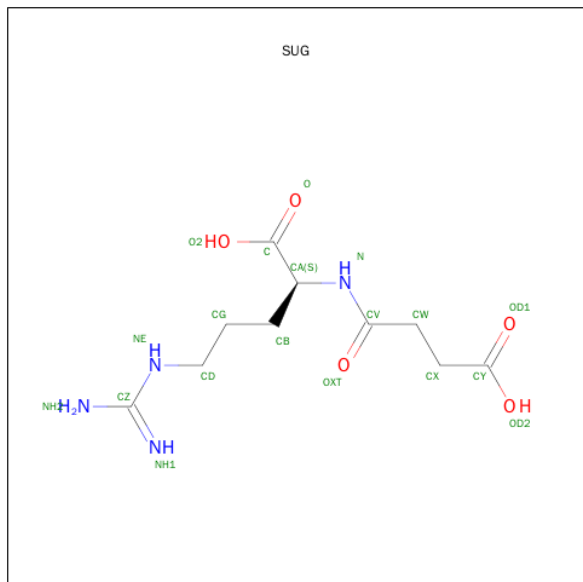
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	HIS	-	EXPRESSION TAG	UNP P76216
B	-1	HIS	-	EXPRESSION TAG	UNP P76216
B	0	GLY	-	EXPRESSION TAG	UNP P76216
B	1	SER	-	EXPRESSION TAG	UNP P76216
B	365	SER	CYS	ENGINEERED	UNP P76216
C	-10	MET	-	EXPRESSION TAG	UNP P76216
C	-9	GLY	-	EXPRESSION TAG	UNP P76216
C	-8	SER	-	EXPRESSION TAG	UNP P76216
C	-7	SER	-	EXPRESSION TAG	UNP P76216
C	-6	HIS	-	EXPRESSION TAG	UNP P76216
C	-5	HIS	-	EXPRESSION TAG	UNP P76216
C	-4	HIS	-	EXPRESSION TAG	UNP P76216
C	-3	HIS	-	EXPRESSION TAG	UNP P76216
C	-2	HIS	-	EXPRESSION TAG	UNP P76216
C	-1	HIS	-	EXPRESSION TAG	UNP P76216
C	0	GLY	-	EXPRESSION TAG	UNP P76216
C	1	SER	-	EXPRESSION TAG	UNP P76216
C	365	SER	CYS	ENGINEERED	UNP P76216
D	-10	MET	-	EXPRESSION TAG	UNP P76216
D	-9	GLY	-	EXPRESSION TAG	UNP P76216
D	-8	SER	-	EXPRESSION TAG	UNP P76216
D	-7	SER	-	EXPRESSION TAG	UNP P76216
D	-6	HIS	-	EXPRESSION TAG	UNP P76216
D	-5	HIS	-	EXPRESSION TAG	UNP P76216
D	-4	HIS	-	EXPRESSION TAG	UNP P76216
D	-3	HIS	-	EXPRESSION TAG	UNP P76216
D	-2	HIS	-	EXPRESSION TAG	UNP P76216
D	-1	HIS	-	EXPRESSION TAG	UNP P76216
D	0	GLY	-	EXPRESSION TAG	UNP P76216
D	1	SER	-	EXPRESSION TAG	UNP P76216
D	365	SER	CYS	ENGINEERED	UNP P76216

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

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

- Molecule 3 is N 2 -(3-CARBOXYPROPANOYL)-L-ARGININE (three-letter code: SUG) (formula: C₁₀H₁₈N₄O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			19	10	4	5		
3	B	1	Total	C	N	O	0	0
			19	10	4	5		
3	C	1	Total	C	N	O	0	0
			19	10	4	5		
3	D	1	Total	C	N	O	0	0
			19	10	4	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	154	Total	O	0	0
			154	154		
4	B	150	Total	O	0	0
			150	150		
4	C	100	Total	O	0	0
			100	100		
4	D	163	Total	O	0	0
			163	163		

3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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: Succinylarginine Dihydrolase

Chain A:

- Molecule 1: Succinylarginine Dihydrolase

Chain B:

71% 22%

0.00 0.05 0.10 0.15

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39 40 41 42 43 44 45 46 47 48 49 50 51 52 53 54 55 56 57 58 59 60 61 62 63 64 65 66 67 68 69 70 71 72 73 74 75 76 77 78 79 80 81 82 83 84 85 86 87 88 89 90 91 92 93 94 95 96 97 98 99 100

Met Gly Ser His His His His His Gly Ser Pz E5 D9 L14 Y18 S22 F23 G24 R32 N37 P38 R39 K43 Q44 G45 M49 K50 D54 V61 H65 F69 I70 P71 V72 L73 S81 Q84 R85 L86 V89 A90 P91

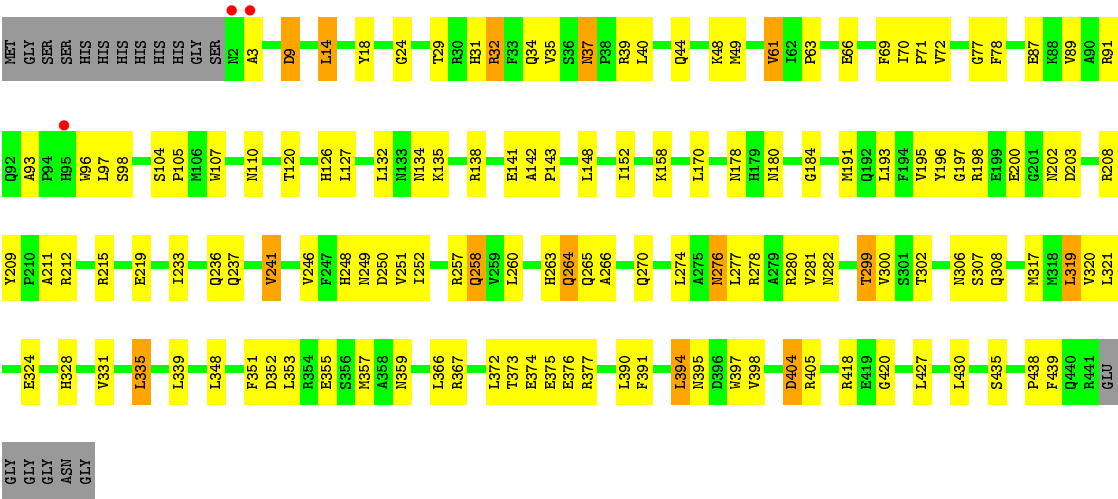
V100 S101 S104 P105 H106 H107 N110 T120 L121 H126 L132 K135 F136 H137 R138 S139 L140 E141 A142 P143 D155 E156 E157 K158 M178 H179 N180 L193 R198 E199 N202 R205 R208 V209 F210 A211 R215 E216 R223 Q226 V227 R228 A229 P230

F234 Q237 V241 V246 H248 N249 D250 V251 I252 R257 Q264 R269 Q270 S271 Q272 L273 L274 A275 N276 E288 Y303 N306 L310 D313 D314 G315 S316 M317 M318 L321 P322 Q323 R326 E327 H328 V331 Y334 L335 E355 S356 M357 P359

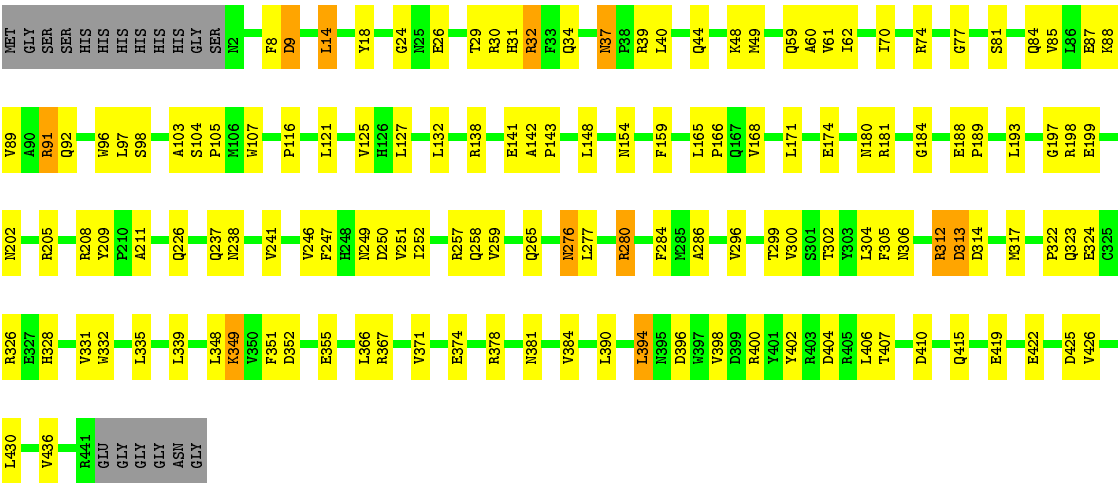
N359 G360 G361 G362 L366 R367 L368 R369 L372 E375 E376 R377 R378 L394 V398 Y402 R405 L411 L430 R441 GLU GLY GLY ASN GLY

- Molecule 1: Succinylarginine Dihydrolase

Chain C:  %



● Molecule 1: Succinylarginine Dihydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.85Å 166.90Å 185.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.74 – 2.20 47.74 – 2.20	Depositor EDS
% Data completeness (in resolution range)	73.9 (47.74-2.20) 73.9 (47.74-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.219 , 0.265 0.218 , 0.265	Depositor DCC
R_{free} test set	3091 reflections (4.77%)	DCC
Wilson B-factor (Å ²)	20.4	Xtriage
Anisotropy	0.318	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 39.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 64862 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14363	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.54 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.2212e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.37	0/3502	0.61	0/4760
1	B	0.37	0/3502	0.61	0/4760
1	C	0.37	0/3502	0.60	0/4760
1	D	0.36	0/3502	0.60	0/4760
All	All	0.37	0/14008	0.60	0/19040

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	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	334	TYR	Sidechain

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	3429	0	3355	100	0
1	B	3429	0	3355	92	0
1	C	3429	0	3355	122	0
1	D	3429	0	3355	107	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	19	0	15	1	0
3	B	19	0	15	4	0
3	C	19	0	15	0	0
3	D	19	0	15	2	0
4	A	154	0	0	3	0
4	B	150	0	0	5	0
4	C	100	0	0	4	0
4	D	163	0	0	7	0
All	All	14363	0	13480	417	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 15.

All (417) 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:A:328:HIS:HD2	1:A:331:VAL:H	1.05	1.04
1:D:14:LEU:H	1:D:14:LEU:HD12	1.23	0.99
1:C:258:GLN:H	1:C:258:GLN:HE21	1.11	0.97
1:C:328:HIS:HD2	1:C:331:VAL:H	1.12	0.97
1:A:241:VAL:HG21	1:A:252:ILE:HD12	1.48	0.95
1:B:328:HIS:HD2	1:B:331:VAL:H	1.07	0.94
1:D:258:GLN:HE21	1:D:259:VAL:HG23	1.36	0.90
1:C:233:ILE:HD11	1:C:280:ARG:HD2	1.54	0.89
1:B:49:MET:HE1	1:B:61:VAL:HG22	1.54	0.89
1:C:258:GLN:NE2	1:C:258:GLN:H	1.69	0.89
1:A:165:LEU:H	1:A:165:LEU:HD12	1.38	0.88
1:C:37:ASN:ND2	1:C:40:LEU:H	1.76	0.84
1:B:198:ARG:NH1	1:B:211:ALA:O	2.11	0.84
1:B:241:VAL:HG21	1:B:252:ILE:CD1	2.08	0.84
1:A:328:HIS:CD2	1:A:331:VAL:H	1.94	0.83
1:A:104:SER:HB2	1:A:138:ARG:HD3	1.60	0.83
1:D:323:GLN:HG3	1:D:352:ASP:HA	1.60	0.83
1:C:104:SER:HB2	1:C:138:ARG:HD3	1.61	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:MET:HE1	1:C:61:VAL:HG12	1.60	0.83
1:D:104:SER:HB2	1:D:138:ARG:HD3	1.61	0.82
1:D:171:LEU:HB2	4:D:2159:HOH:O	1.79	0.82
1:B:328:HIS:CD2	1:B:331:VAL:H	1.97	0.81
1:D:328:HIS:HD2	1:D:331:VAL:H	1.25	0.81
1:B:49:MET:CE	1:B:61:VAL:HG22	2.09	0.81
1:A:241:VAL:HG21	1:A:252:ILE:CD1	2.12	0.79
1:B:14:LEU:HD12	1:B:14:LEU:H	1.47	0.79
1:D:14:LEU:N	1:D:14:LEU:HD12	1.99	0.78
1:A:138:ARG:HD2	1:A:141:GLU:OE2	1.83	0.78
1:B:357:MET:HE2	1:B:357:MET:HA	1.66	0.78
1:C:252:ILE:HD13	1:C:266:ALA:HA	1.66	0.76
1:B:155:ASP:OD1	1:B:157:GLU:HG3	1.86	0.75
1:C:215:ARG:O	1:C:219:GLU:HG3	1.87	0.74
1:C:258:GLN:HE21	1:C:258:GLN:N	1.85	0.73
1:D:88:LYS:HB3	1:D:92:GLN:NE2	2.04	0.72
1:A:165:LEU:N	1:A:165:LEU:HD12	2.05	0.71
1:C:110:ASN:HD21	1:C:248:HIS:CE1	2.08	0.71
1:D:14:LEU:H	1:D:14:LEU:CD1	1.99	0.71
1:D:238:ASN:O	1:D:241:VAL:HG12	1.90	0.71
1:C:39:ARG:NH2	1:C:404:ASP:HB3	2.07	0.70
1:D:168:VAL:HG22	4:D:2159:HOH:O	1.90	0.69
1:D:142:ALA:HB3	1:D:143:PRO:HD3	1.74	0.69
1:A:178:ASN:ND2	1:A:249:ASN:H	1.89	0.69
1:B:228:ASN:HD22	1:B:229:PRO:CD	2.06	0.69
1:C:35:VAL:HG11	1:C:405:ARG:HG3	1.75	0.69
1:A:338:LEU:HD11	1:A:345:ILE:HD12	1.74	0.69
1:C:138:ARG:HD2	1:C:141:GLU:OE2	1.93	0.68
1:C:14:LEU:HD12	1:C:14:LEU:H	1.58	0.68
1:B:314:ASP:OD1	1:B:316:SER:HB3	1.93	0.68
1:A:84:GLN:O	1:A:88:LYS:HG3	1.94	0.68
1:B:394:LEU:O	1:B:398:VAL:HG23	1.93	0.68
1:B:228:ASN:HD22	1:B:229:PRO:HD2	1.59	0.67
1:B:142:ALA:HB3	1:B:143:PRO:HD3	1.77	0.67
1:A:165:LEU:H	1:A:165:LEU:CD1	2.07	0.67
1:D:138:ARG:HD2	1:D:141:GLU:OE2	1.95	0.67
1:D:302:THR:HG21	1:D:331:VAL:HG21	1.77	0.67
1:D:394:LEU:O	1:D:398:VAL:HG23	1.95	0.67
1:C:142:ALA:HB3	1:C:143:PRO:HD3	1.74	0.67
1:A:338:LEU:HD21	1:A:345:ILE:HD13	1.75	0.67
1:A:220:ALA:HA	1:A:223:ARG:NH1	2.10	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:241:VAL:HG13	1:C:246:VAL:HG21	1.76	0.66
1:D:237:GLN:HE22	1:D:249:ASN:HD22	1.43	0.66
1:C:237:GLN:HE22	1:C:249:ASN:ND2	1.94	0.66
1:C:328:HIS:CD2	1:C:331:VAL:HG23	2.31	0.65
1:C:241:VAL:HG13	1:C:246:VAL:CG2	2.27	0.65
1:D:390:LEU:HG	1:D:394:LEU:HD22	1.77	0.65
1:A:49:MET:CE	1:A:61:VAL:HG12	2.27	0.65
1:C:319:LEU:HD13	1:C:321:LEU:HD21	1.78	0.65
1:A:49:MET:HE1	1:A:61:VAL:HG12	1.80	0.64
1:D:81:SER:OG	1:D:84:GLN:HG3	1.97	0.64
1:B:216:GLU:N	1:B:216:GLU:OE2	2.24	0.64
1:D:258:GLN:NE2	1:D:259:VAL:HG23	2.10	0.63
1:B:180:ASN:HB2	1:B:193:LEU:HB3	1.79	0.63
1:D:29:THR:O	1:D:32:ARG:HG3	1.97	0.63
1:D:37:ASN:ND2	1:D:40:LEU:H	1.95	0.63
1:C:233:ILE:HD11	1:C:280:ARG:CD	2.28	0.62
1:C:69:PHE:O	1:C:72:VAL:HG12	1.99	0.62
1:A:338:LEU:HD21	1:A:345:ILE:CD1	2.29	0.62
1:D:324:GLU:H	1:D:324:GLU:CD	2.01	0.62
1:C:35:VAL:CG1	1:C:405:ARG:HG3	2.29	0.62
1:D:84:GLN:O	1:D:88:LYS:HG3	2.00	0.62
1:C:241:VAL:HG21	1:C:252:ILE:CD1	2.30	0.62
1:B:14:LEU:HD12	1:B:14:LEU:N	2.15	0.61
1:C:307:SER:HB3	1:C:321:LEU:HD23	1.82	0.61
1:A:142:ALA:HB3	1:A:143:PRO:HD3	1.82	0.61
1:C:328:HIS:CD2	1:C:331:VAL:H	2.05	0.61
1:A:50:LYS:HE3	1:A:54:ASP:OD2	2.01	0.61
1:D:14:LEU:HG	1:D:402:TYR:CE1	2.36	0.60
1:D:349:LYS:HG2	1:D:351:PHE:CE1	2.36	0.60
1:A:27:ALA:O	1:A:31:HIS:HD2	1.83	0.60
1:C:237:GLN:HE22	1:C:249:ASN:HD22	1.46	0.60
1:B:110:ASN:HD21	1:B:248:HIS:CE1	2.20	0.60
1:B:14:LEU:CD1	1:B:14:LEU:H	2.14	0.60
1:C:331:VAL:HG12	1:C:335:LEU:HD22	1.83	0.60
1:B:241:VAL:HG21	1:B:252:ILE:HD13	1.83	0.60
1:A:44:GLN:HE22	1:A:355:GLU:H	1.47	0.60
1:B:357:MET:HE1	4:B:2096:HOH:O	2.02	0.59
1:C:104:SER:CB	1:C:138:ARG:HD3	2.33	0.59
1:C:274:LEU:O	1:C:278:ARG:HG3	2.02	0.59
1:C:158:LYS:HE2	1:C:375:GLU:HG2	1.85	0.59
1:A:104:SER:CB	1:A:138:ARG:HD3	2.32	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:44:GLN:HE22	1:C:355:GLU:H	1.51	0.59
1:A:169:ALA:O	1:A:198:ARG:NH2	2.33	0.59
1:A:280:ARG:HH11	1:A:280:ARG:HG3	1.67	0.59
1:B:138:ARG:HD3	1:B:141:GLU:OE2	2.03	0.59
1:A:166:PRO:HD2	1:A:171:LEU:HD13	1.83	0.58
1:D:241:VAL:HG22	1:D:246:VAL:HG21	1.85	0.58
1:C:191:MET:SD	1:C:280:ARG:HD3	2.44	0.58
1:D:322:PRO:HB2	1:D:324:GLU:OE2	2.04	0.58
1:A:204:THR:HG22	1:A:236:GLN:HE22	1.69	0.58
1:D:44:GLN:HE22	1:D:355:GLU:H	1.50	0.57
1:B:69:PHE:CD2	1:B:72:VAL:HG23	2.40	0.57
1:C:14:LEU:H	1:C:14:LEU:CD1	2.17	0.57
1:C:93:ALA:HB1	1:C:96:TRP:HD1	1.69	0.57
1:A:237:GLN:HE22	1:A:249:ASN:HD22	1.51	0.57
1:D:104:SER:HA	1:D:107:TRP:CZ3	2.40	0.56
1:B:69:PHE:HD2	1:B:72:VAL:HG23	1.70	0.56
1:B:81:SER:OG	1:B:84:GLN:HG3	2.05	0.56
1:C:110:ASN:HB3	1:C:132:LEU:HD21	1.86	0.56
1:D:415:GLN:HG3	1:D:419:GLU:OE1	2.05	0.56
1:B:228:ASN:HD22	1:B:229:PRO:N	2.02	0.56
1:A:39:ARG:HD2	1:A:402:TYR:O	2.05	0.56
1:D:88:LYS:HB3	1:D:92:GLN:HE21	1.70	0.56
1:D:328:HIS:CD2	1:D:331:VAL:H	2.16	0.56
1:C:404:ASP:OD2	1:C:404:ASP:N	2.28	0.56
1:A:289:VAL:HG22	1:A:334:TYR:CZ	2.41	0.56
1:D:39:ARG:HG3	1:D:398:VAL:HG12	1.88	0.56
1:C:418:ARG:HH11	1:C:418:ARG:HG3	1.70	0.56
1:D:202:ASN:HB3	4:D:2066:HOH:O	2.05	0.56
1:C:31:HIS:ND1	1:C:34:GLN:NE2	2.53	0.56
1:B:178:ASN:ND2	1:B:249:ASN:H	2.04	0.56
1:A:308:GLN:HB2	1:A:320:VAL:HB	1.86	0.55
1:A:372:LEU:HB2	1:A:377:ARG:HG3	1.87	0.55
1:A:25:ASN:HD22	1:A:25:ASN:C	2.10	0.55
1:B:223:ARG:HG2	1:C:77:GLY:O	2.06	0.55
1:A:49:MET:HE1	1:A:61:VAL:CG1	2.37	0.55
1:C:324:GLU:HG3	1:C:357:MET:CE	2.37	0.55
1:D:166:PRO:O	4:D:2159:HOH:O	2.18	0.55
1:C:110:ASN:ND2	1:C:248:HIS:CE1	2.74	0.55
1:B:228:ASN:ND2	1:B:229:PRO:HD2	2.22	0.55
1:A:39:ARG:HD3	1:A:404:ASP:OD1	2.07	0.55
1:B:22:SER:HA	1:B:135:LYS:HE3	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:ARG:HG3	1:A:92:GLN:HG3	1.89	0.54
1:A:127:LEU:N	1:A:127:LEU:HD12	2.23	0.54
1:B:269:ARG:HB3	1:B:272:GLN:HG2	1.89	0.54
1:A:224:LEU:HD23	1:D:77:GLY:HA2	1.89	0.54
1:D:323:GLN:NE2	1:D:352:ASP:HB2	2.22	0.54
1:C:324:GLU:CD	1:C:324:GLU:H	2.11	0.54
1:C:14:LEU:N	1:C:14:LEU:HD12	2.21	0.54
1:C:397:TRP:CH2	1:C:427:LEU:HD22	2.43	0.54
1:D:326:ARG:HD2	4:D:2062:HOH:O	2.08	0.54
1:C:373:THR:OG1	1:C:376:GLU:HG3	2.08	0.54
1:B:272:GLN:HG3	1:B:273:LEU:N	2.22	0.54
1:D:85:VAL:O	1:D:89:VAL:HG23	2.08	0.53
1:D:339:LEU:HD22	1:D:348:LEU:HG	1.91	0.53
1:C:352:ASP:C	1:C:353:LEU:HD12	2.28	0.53
1:A:280:ARG:NH1	1:A:280:ARG:HG3	2.24	0.53
1:C:37:ASN:C	1:C:37:ASN:HD22	2.12	0.53
1:A:400:ARG:NE	1:A:401:TYR:CE1	2.77	0.53
1:A:321:LEU:HD13	1:A:335:LEU:HD23	1.91	0.53
1:A:104:SER:HA	1:A:107:TRP:CZ3	2.44	0.52
1:B:138:ARG:HH21	3:B:1002:SUG:HG2	1.75	0.52
1:A:241:VAL:HG22	1:A:246:VAL:HG21	1.91	0.52
1:B:251:VAL:O	1:B:251:VAL:HG12	2.08	0.52
1:C:241:VAL:HG21	1:C:252:ILE:HD12	1.92	0.52
1:B:246:VAL:HG13	1:B:251:VAL:HB	1.91	0.52
1:B:237:GLN:HE22	1:B:249:ASN:HD22	1.57	0.52
1:C:339:LEU:HD22	1:C:348:LEU:HG	1.92	0.52
1:C:237:GLN:NE2	1:C:249:ASN:HD22	2.08	0.52
1:C:251:VAL:HG12	1:C:251:VAL:O	2.09	0.52
1:A:191:MET:HE1	1:A:233:ILE:HG13	1.91	0.52
1:B:158:LYS:HE3	1:B:375:GLU:OE2	2.10	0.52
1:D:252:ILE:HD12	1:D:299:THR:HG21	1.92	0.52
1:D:331:VAL:HG12	1:D:335:LEU:HD23	1.92	0.51
1:A:126:HIS:C	1:A:127:LEU:HD12	2.31	0.51
1:C:93:ALA:HB1	1:C:96:TRP:CD1	2.45	0.51
1:C:260:LEU:C	1:C:260:LEU:HD23	2.30	0.51
1:C:331:VAL:HG12	1:C:335:LEU:CD2	2.40	0.51
1:D:251:VAL:O	1:D:251:VAL:HG12	2.11	0.51
1:B:101:SER:HB3	1:B:411:LEU:HD22	1.92	0.51
1:C:24:GLY:HA2	1:C:209:TYR:CD2	2.46	0.51
1:B:326:ARG:HD2	1:B:326:ARG:O	2.11	0.51
1:A:96:TRP:CE3	1:A:96:TRP:HA	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:ASN:HD22	1:B:37:ASN:C	2.13	0.51
1:D:265:GLN:HA	4:D:2065:HOH:O	2.10	0.51
1:A:328:HIS:HD2	1:A:331:VAL:N	1.89	0.50
1:D:104:SER:CB	1:D:138:ARG:HD3	2.36	0.50
1:A:180:ASN:HB2	1:A:193:LEU:HB3	1.93	0.50
1:C:264:GLN:HG3	1:C:265:GLN:HG3	1.92	0.50
1:D:49:MET:HE2	1:D:59:GLN:HE21	1.76	0.50
1:D:415:GLN:O	1:D:419:GLU:HG3	2.11	0.50
1:A:18:TYR:CE2	1:A:358:ALA:HB3	2.46	0.50
1:A:9:ASP:HA	1:A:366:LEU:O	2.12	0.50
1:A:238:ASN:O	1:A:241:VAL:HG12	2.11	0.50
1:B:228:ASN:C	1:B:228:ASN:HD22	2.15	0.50
1:A:252:ILE:HD13	1:A:266:ALA:HA	1.93	0.49
1:A:104:SER:N	1:A:105:PRO:HD2	2.26	0.49
1:D:331:VAL:O	1:D:335:LEU:HD23	2.12	0.49
1:C:48:LYS:HA	1:C:351:PHE:CE2	2.47	0.49
1:C:120:THR:HG21	1:C:126:HIS:HD2	1.76	0.49
1:D:39:ARG:HD3	1:D:404:ASP:OD2	2.12	0.49
1:A:390:LEU:HG	1:A:394:LEU:HD22	1.94	0.49
1:D:116:PRO:HG3	1:D:181:ARG:CZ	2.41	0.49
1:C:9:ASP:HA	1:C:366:LEU:O	2.12	0.49
1:A:343:ASN:HB2	1:A:344:PRO:CD	2.42	0.49
1:B:49:MET:HE3	1:B:61:VAL:HG22	1.92	0.49
1:B:257:ARG:NH2	4:B:2065:HOH:O	2.46	0.49
1:A:8:PHE:CD2	1:A:60:ALA:HB3	2.47	0.49
1:A:110:ASN:HD21	1:A:248:HIS:CE1	2.30	0.49
1:D:104:SER:N	1:D:105:PRO:HD2	2.27	0.49
1:C:241:VAL:HG13	1:C:246:VAL:CB	2.42	0.49
1:C:418:ARG:NH1	1:C:418:ARG:HG3	2.27	0.49
1:B:110:ASN:ND2	1:B:248:HIS:CE1	2.80	0.49
1:B:241:VAL:HG22	1:B:246:VAL:HG21	1.94	0.49
1:D:44:GLN:NE2	4:D:2091:HOH:O	2.46	0.49
1:D:9:ASP:HA	1:D:366:LEU:O	2.12	0.49
1:C:148:LEU:O	1:C:152:ILE:HD13	2.13	0.48
1:A:269:ARG:NH1	1:A:272:GLN:HE22	2.11	0.48
1:B:50:LYS:NZ	1:B:54:ASP:OD1	2.47	0.48
1:C:78:PHE:CD1	1:C:89:VAL:HG22	2.48	0.48
1:A:197:GLY:HA3	1:A:211:ALA:HB1	1.95	0.48
1:B:44:GLN:HE22	1:B:355:GLU:H	1.62	0.48
1:B:9:ASP:HA	1:B:366:LEU:O	2.13	0.48
1:B:264:GLN:HB2	1:B:288:GLU:OE2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:37:ASN:C	1:D:37:ASN:HD22	2.18	0.47
1:B:14:LEU:HG	1:B:402:TYR:CE1	2.49	0.47
1:D:241:VAL:HG22	1:D:246:VAL:CG2	2.44	0.47
1:A:227:VAL:HG11	1:A:232:VAL:HG22	1.95	0.47
1:D:96:TRP:CE3	1:D:96:TRP:HA	2.49	0.47
1:C:191:MET:CE	1:C:280:ARG:HG2	2.44	0.47
1:A:237:GLN:HE22	1:A:249:ASN:ND2	2.12	0.47
1:C:14:LEU:HD22	1:C:420:GLY:HA2	1.96	0.47
1:D:188:GLU:HB3	1:D:189:PRO:HD2	1.97	0.47
1:D:258:GLN:HE22	1:D:317:MET:CE	2.28	0.47
1:C:104:SER:N	1:C:105:PRO:HD2	2.29	0.47
1:C:241:VAL:HG13	1:C:246:VAL:HB	1.96	0.47
1:A:121:LEU:HB2	1:A:226:GLN:HE21	1.79	0.47
1:C:91:ARG:HH11	1:C:91:ARG:HB3	1.80	0.47
1:D:26:GLU:O	1:D:30:ARG:HG3	2.15	0.47
1:D:284:PHE:CE1	1:D:286:ALA:HB2	2.49	0.47
1:C:282:ASN:N	4:C:2094:HOH:O	2.48	0.47
1:B:215:ARG:HH12	1:B:234:PHE:HB2	1.80	0.47
1:B:121:LEU:HD12	1:B:226:GLN:HG2	1.96	0.47
1:B:303:TYR:CD2	1:B:357:MET:HE3	2.49	0.47
1:B:270:GLN:O	1:B:274:LEU:HG	2.15	0.47
1:C:208:ARG:HH11	1:C:208:ARG:HG2	1.79	0.46
1:A:83:GLU:O	1:A:87:GLU:HG3	2.15	0.46
1:B:107:TRP:CD2	3:B:1002:SUG:HG1	2.50	0.46
1:C:263:HIS:ND1	1:C:264:GLN:N	2.63	0.46
1:D:180:ASN:HB2	1:D:193:LEU:HB3	1.97	0.46
1:A:14:LEU:HD21	1:A:420:GLY:HA2	1.97	0.46
1:C:107:TRP:CH2	1:C:359:ASN:HB3	2.50	0.46
1:B:199:GLU:HG3	1:B:202:ASN:HB2	1.97	0.46
1:C:390:LEU:HG	1:C:394:LEU:HD22	1.97	0.46
1:C:37:ASN:HD22	1:C:40:LEU:H	1.61	0.46
1:B:105:PRO:HB3	1:B:141:GLU:HG2	1.97	0.46
1:A:66:GLU:CD	1:A:439:PHE:HB3	2.35	0.46
1:B:223:ARG:HB3	1:C:77:GLY:HA3	1.98	0.46
1:A:18:TYR:O	1:A:32:ARG:O	2.34	0.46
1:A:331:VAL:O	1:A:335:LEU:HD13	2.16	0.46
1:C:178:ASN:ND2	1:C:249:ASN:H	2.13	0.46
1:C:257:ARG:HB2	1:C:317:MET:HE1	1.98	0.45
1:B:357:MET:HE3	1:B:362:GLY:CA	2.46	0.45
1:C:264:GLN:HA	1:C:270:GLN:HE22	1.80	0.45
1:D:247:PHE:HE1	3:D:1004:SUG:HD1	1.79	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:432:ASN:ND2	4:A:2014:HOH:O	2.48	0.45
1:C:299:THR:HG22	1:C:300:VAL:N	2.31	0.45
1:C:302:THR:HG21	1:C:331:VAL:HG21	1.98	0.45
1:B:104:SER:N	1:B:105:PRO:HD2	2.31	0.45
1:C:104:SER:HA	1:C:107:TRP:CZ3	2.51	0.45
1:C:39:ARG:HG3	1:C:398:VAL:HG12	1.97	0.45
1:C:29:THR:O	1:C:32:ARG:HG3	2.17	0.45
1:C:331:VAL:O	1:C:335:LEU:HD22	2.16	0.45
1:C:49:MET:CE	1:C:61:VAL:HG12	2.38	0.45
1:B:107:TRP:CH2	1:B:359:ASN:HB3	2.52	0.45
1:B:24:GLY:HA2	1:B:209:TYR:CG	2.52	0.45
1:D:332:TRP:CE3	1:D:332:TRP:HA	2.52	0.45
1:B:357:MET:HE2	1:B:361:GLY:O	2.16	0.45
1:A:81:SER:OG	1:A:84:GLN:HG3	2.16	0.45
1:C:324:GLU:HG3	1:C:357:MET:HE1	1.99	0.45
1:B:73:LEU:HD21	1:B:100:VAL:HG21	1.99	0.44
1:B:241:VAL:HG21	1:B:252:ILE:HD12	1.97	0.44
1:D:31:HIS:HA	1:D:34:GLN:NE2	2.31	0.44
1:D:70:ILE:HG22	1:D:74:ARG:NH1	2.32	0.44
1:D:296:VAL:O	1:D:300:VAL:HG23	2.17	0.44
1:D:422:GLU:O	1:D:426:VAL:HG23	2.17	0.44
1:B:39:ARG:O	1:B:43:LYS:HG2	2.17	0.44
1:A:153:PHE:HB3	1:A:159:PHE:CG	2.52	0.44
1:C:184:GLY:O	1:C:257:ARG:HD2	2.18	0.44
1:A:241:VAL:CG2	1:A:252:ILE:HD12	2.32	0.44
1:D:349:LYS:HG2	1:D:351:PHE:HE1	1.80	0.44
1:B:72:VAL:HG21	1:B:140:LEU:HD22	2.00	0.44
1:C:18:TYR:O	1:C:32:ARG:O	2.35	0.44
1:C:257:ARG:HB3	1:C:258:GLN:NE2	2.32	0.44
1:A:394:LEU:O	1:A:398:VAL:HG23	2.17	0.44
1:C:91:ARG:CB	1:C:91:ARG:NH1	2.80	0.44
1:D:8:PHE:CD2	1:D:60:ALA:HB3	2.52	0.44
1:D:166:PRO:HD2	1:D:171:LEU:HD13	1.99	0.44
1:A:191:MET:HE1	1:A:232:VAL:C	2.38	0.44
1:D:127:LEU:N	1:D:127:LEU:HD12	2.33	0.44
1:B:303:TYR:HA	4:B:2004:HOH:O	2.17	0.44
1:A:96:TRP:HE3	1:A:96:TRP:HA	1.82	0.44
1:C:197:GLY:HA3	1:C:211:ALA:HB1	2.00	0.44
1:D:184:GLY:O	1:D:257:ARG:HD2	2.17	0.44
1:C:66:GLU:OE1	1:C:438:PRO:HD2	2.18	0.44
1:C:93:ALA:CB	1:C:96:TRP:HD1	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:ALA:O	1:A:31:HIS:CD2	2.68	0.44
1:C:170:LEU:HD11	1:C:200:GLU:CD	2.38	0.44
1:B:120:THR:HG21	1:B:126:HIS:HD2	1.83	0.43
1:C:195:VAL:HG11	1:C:249:ASN:HD22	1.83	0.43
1:A:158:LYS:HE2	1:A:375:GLU:HG2	1.99	0.43
1:A:107:TRP:CD2	3:A:1001:SUG:HG1	2.53	0.43
1:A:204:THR:HG22	1:A:236:GLN:NE2	2.34	0.43
1:D:323:GLN:HG3	1:D:352:ASP:CA	2.39	0.43
1:D:198:ARG:NH1	1:D:211:ALA:O	2.51	0.43
1:A:191:MET:HE3	1:A:192:GLN:N	2.34	0.43
1:C:281:VAL:HG13	4:C:2094:HOH:O	2.18	0.43
1:B:44:GLN:NE2	4:B:2019:HOH:O	2.48	0.43
1:B:198:ARG:HG2	1:B:198:ARG:HH11	1.83	0.43
1:B:198:ARG:HD3	1:B:205:ARG:NH2	2.34	0.43
1:C:66:GLU:OE2	1:C:439:PHE:HB3	2.19	0.43
1:A:299:THR:HG23	1:A:305:PHE:HD1	1.84	0.43
1:D:132:LEU:HD23	1:D:174:GLU:HA	2.00	0.43
1:C:250:ASP:O	1:C:306:ASN:HA	2.18	0.43
1:B:367:ARG:HD2	1:B:367:ARG:C	2.39	0.43
1:C:252:ILE:HG12	1:C:252:ILE:O	2.19	0.43
1:A:224:LEU:CD2	1:D:77:GLY:HA2	2.49	0.43
1:B:208:ARG:HD3	1:B:209:TYR:CE2	2.54	0.43
1:A:70:ILE:N	1:A:71:PRO:CD	2.82	0.43
1:A:125:VAL:HB	1:A:159:PHE:HD2	1.84	0.43
1:D:312:ARG:HG3	1:D:312:ARG:HH11	1.84	0.43
1:D:24:GLY:HA2	1:D:209:TYR:CG	2.54	0.43
1:C:308:GLN:HB2	1:C:320:VAL:HB	2.00	0.42
1:B:107:TRP:HB2	1:B:138:ARG:HH22	1.83	0.42
1:C:31:HIS:O	1:C:32:ARG:C	2.58	0.42
1:D:312:ARG:CG	1:D:312:ARG:HH11	2.32	0.42
1:C:257:ARG:HH11	1:C:257:ARG:HG3	1.84	0.42
1:B:247:PHE:HE1	3:B:1002:SUG:HD1	1.83	0.42
1:A:190:GLY:O	1:A:231:GLN:HB3	2.19	0.42
1:B:85:VAL:O	1:B:89:VAL:HG23	2.18	0.42
1:D:276:ASN:O	1:D:280:ARG:CG	2.67	0.42
1:C:202:ASN:HB3	4:C:2049:HOH:O	2.18	0.42
1:C:70:ILE:N	1:C:71:PRO:CD	2.82	0.42
1:C:134:ASN:OD1	1:C:135:LYS:HE2	2.20	0.42
1:D:18:TYR:O	1:D:32:ARG:O	2.38	0.42
1:D:125:VAL:HB	1:D:159:PHE:HD2	1.84	0.42
1:D:406:LEU:HD12	1:D:410:ASP:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:212:ARG:HB2	4:C:2063:HOH:O	2.19	0.42
1:D:103:ALA:C	1:D:105:PRO:HD2	2.40	0.42
1:C:324:GLU:HG3	1:C:357:MET:HE2	2.02	0.42
1:D:184:GLY:O	1:D:257:ARG:CD	2.68	0.42
1:A:70:ILE:HG12	4:A:2017:HOH:O	2.20	0.42
1:D:121:LEU:HB2	1:D:226:GLN:HE21	1.84	0.42
1:A:46:LEU:HD23	1:A:49:MET:HE2	2.02	0.42
1:B:136:PHE:CE2	1:B:140:LEU:HD21	2.55	0.42
1:D:198:ARG:HB2	1:D:205:ARG:HH11	1.84	0.42
1:A:250:ASP:O	1:A:306:ASN:HA	2.20	0.42
1:C:196:TYR:CE2	1:C:215:ARG:HD2	2.55	0.42
1:B:5:GLU:OE2	1:B:369:ARG:HD2	2.19	0.42
1:D:314:ASP:C	1:D:314:ASP:OD2	2.59	0.42
1:A:403:ARG:HG3	4:A:2089:HOH:O	2.20	0.42
1:C:63:PRO:CG	1:C:148:LEU:HD22	2.49	0.41
1:D:331:VAL:HG12	1:D:335:LEU:CD2	2.50	0.41
1:D:246:VAL:HG13	1:D:251:VAL:HB	2.02	0.41
1:A:25:ASN:ND2	1:A:28:SER:H	2.18	0.41
1:C:264:GLN:HB2	1:C:264:GLN:HE21	1.59	0.41
1:D:116:PRO:HB3	1:D:371:VAL:HB	2.02	0.41
1:B:372:LEU:HB2	1:B:377:ARG:HG2	2.01	0.41
1:C:3:ALA:HB1	1:C:372:LEU:O	2.20	0.41
1:A:107:TRP:HD1	1:A:365:SER:HG	1.68	0.41
1:B:378:ARG:HA	1:B:378:ARG:NE	2.35	0.41
1:C:276:ASN:ND2	1:C:276:ASN:C	2.73	0.41
1:C:374:GLU:HA	1:C:377:ARG:CZ	2.50	0.41
1:D:396:ASP:O	1:D:400:ARG:HG3	2.19	0.41
1:C:258:GLN:NE2	1:C:258:GLN:N	2.51	0.41
1:A:105:PRO:HA	1:A:141:GLU:OE1	2.20	0.41
1:C:195:VAL:HG11	1:C:249:ASN:ND2	2.35	0.41
1:B:132:LEU:HD13	1:B:138:ARG:HD2	2.03	0.41
1:B:250:ASP:O	1:B:306:ASN:HA	2.20	0.41
1:A:251:VAL:HG12	1:A:251:VAL:O	2.20	0.41
1:A:324:GLU:HG3	1:A:357:MET:CE	2.50	0.41
1:B:49:MET:HE2	1:B:49:MET:HB3	1.84	0.41
1:D:247:PHE:CE1	3:D:1004:SUG:HD1	2.54	0.41
1:D:148:LEU:HD23	1:D:436:VAL:HG21	2.02	0.41
1:A:310:LEU:HD12	1:A:318:MET:HE2	2.02	0.41
1:C:391:PHE:O	1:C:395:ASN:ND2	2.53	0.41
1:C:127:LEU:HD12	1:C:127:LEU:N	2.36	0.41
1:D:14:LEU:HG	1:D:402:TYR:CZ	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:328:HIS:HD2	1:B:331:VAL:N	1.92	0.41
1:D:48:LYS:HA	1:D:351:PHE:CE2	2.55	0.41
1:B:18:TYR:O	1:B:32:ARG:O	2.38	0.41
1:B:70:ILE:HD11	1:B:86:LEU:HD13	2.03	0.41
1:C:307:SER:CB	1:C:321:LEU:HD23	2.48	0.41
1:D:199:GLU:HG3	1:D:202:ASN:HB2	2.03	0.41
1:D:381:ASN:O	1:D:384:VAL:HG22	2.21	0.41
1:B:247:PHE:CE1	3:B:1002:SUG:HD1	2.56	0.41
1:D:312:ARG:HD3	1:D:313:ASP:H	1.86	0.41
1:B:45:GLY:HA3	4:B:2024:HOH:O	2.20	0.41
1:D:250:ASP:O	1:D:306:ASN:HA	2.19	0.41
1:B:335:LEU:HA	1:B:335:LEU:HD13	1.93	0.41
1:A:103:ALA:C	1:A:105:PRO:HD2	2.41	0.41
1:A:19:ALA:HB3	1:A:21:LEU:HG	2.03	0.41
1:C:180:ASN:HB2	1:C:193:LEU:HB3	2.02	0.41
1:A:394:LEU:HA	1:A:394:LEU:HD12	1.85	0.40
1:D:96:TRP:HA	1:D:96:TRP:HE3	1.86	0.40
1:A:241:VAL:HG22	1:A:246:VAL:CG2	2.52	0.40
1:B:328:HIS:CD2	1:B:331:VAL:HG23	2.56	0.40
1:D:138:ARG:HA	1:D:141:GLU:OE2	2.21	0.40
1:D:197:GLY:HA3	1:D:211:ALA:HB1	2.02	0.40
1:A:199:GLU:HG3	1:A:202:ASN:HB2	2.03	0.40
1:B:310:LEU:HD12	1:B:318:MET:HE2	2.03	0.40
1:A:220:ALA:HA	1:A:223:ARG:HH11	1.84	0.40
1:D:31:HIS:O	1:D:32:ARG:C	2.60	0.40
1:C:236:GLN:HG2	1:C:237:GLN:O	2.21	0.40
1:D:61:VAL:CG2	1:D:62:ILE:N	2.84	0.40
1:D:299:THR:HG23	1:D:305:PHE:HD1	1.85	0.40
1:A:206:PRO:HG3	1:A:211:ALA:HB2	2.03	0.40
1:D:87:GLU:O	1:D:91:ARG:HG3	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	438/458 (96%)	422 (96%)	15 (3%)	1 (0%)	52	59
1	B	438/458 (96%)	418 (95%)	19 (4%)	1 (0%)	52	59
1	C	438/458 (96%)	420 (96%)	17 (4%)	1 (0%)	52	59
1	D	438/458 (96%)	420 (96%)	17 (4%)	1 (0%)	52	59
All	All	1752/1832 (96%)	1680 (96%)	68 (4%)	4 (0%)	52	59

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	32	ARG
1	B	32	ARG
1	C	32	ARG
1	D	32	ARG

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	364/377 (97%)	354 (97%)	10 (3%)	52	64
1	B	364/377 (97%)	344 (94%)	20 (6%)	27	30
1	C	364/377 (97%)	342 (94%)	22 (6%)	24	26
1	D	364/377 (97%)	341 (94%)	23 (6%)	22	24
All	All	1456/1508 (97%)	1381 (95%)	75 (5%)	29	33

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ASP
1	A	25	ASN
1	A	43	LYS
1	A	96	TRP

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Mol	Chain	Res	Type
1	A	277	LEU
1	A	312	ARG
1	A	313	ASP
1	A	367	ARG
1	A	394	LEU
1	A	429	GLN
1	B	9	ASP
1	B	14	LEU
1	B	37	ASN
1	B	65	HIS
1	B	91	ARG
1	B	157	GLU
1	B	208	ARG
1	B	226	GLN
1	B	228	ASN
1	B	241	VAL
1	B	264	GLN
1	B	272	GLN
1	B	276	ASN
1	B	316	SER
1	B	321	LEU
1	B	323	GLN
1	B	367	ARG
1	B	378	ARG
1	B	405	ARG
1	B	430	LEU
1	C	9	ASP
1	C	14	LEU
1	C	37	ASN
1	C	61	VAL
1	C	87	GLU
1	C	97	LEU
1	C	98	SER
1	C	198	ARG
1	C	203	ASP
1	C	241	VAL
1	C	258	GLN
1	C	264	GLN
1	C	276	ASN
1	C	277	LEU
1	C	299	THR
1	C	319	LEU

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Mol	Chain	Res	Type
1	C	335	LEU
1	C	367	ARG
1	C	394	LEU
1	C	404	ASP
1	C	430	LEU
1	C	435	SER
1	D	9	ASP
1	D	14	LEU
1	D	37	ASN
1	D	91	ARG
1	D	97	LEU
1	D	98	SER
1	D	154	ASN
1	D	165	LEU
1	D	208	ARG
1	D	276	ASN
1	D	277	LEU
1	D	280	ARG
1	D	304	LEU
1	D	312	ARG
1	D	313	ASP
1	D	349	LYS
1	D	367	ARG
1	D	374	GLU
1	D	378	ARG
1	D	394	LEU
1	D	407	THR
1	D	425	ASP
1	D	430	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (50) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	44	GLN
1	A	178	ASN
1	A	226	GLN
1	A	236	GLN
1	A	249	ASN
1	A	272	GLN
1	A	276	ASN
1	A	328	HIS

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Mol	Chain	Res	Type
1	A	432	ASN
1	B	37	ASN
1	B	44	GLN
1	B	65	HIS
1	B	84	GLN
1	B	178	ASN
1	B	185	HIS
1	B	228	ASN
1	B	230	GLN
1	B	236	GLN
1	B	249	ASN
1	B	264	GLN
1	B	272	GLN
1	B	276	ASN
1	B	323	GLN
1	B	328	HIS
1	C	34	GLN
1	C	37	ASN
1	C	44	GLN
1	C	84	GLN
1	C	178	ASN
1	C	226	GLN
1	C	248	HIS
1	C	249	ASN
1	C	258	GLN
1	C	264	GLN
1	C	276	ASN
1	C	328	HIS
1	D	37	ASN
1	D	44	GLN
1	D	59	GLN
1	D	84	GLN
1	D	92	GLN
1	D	154	ASN
1	D	178	ASN
1	D	226	GLN
1	D	249	ASN
1	D	258	GLN
1	D	276	ASN
1	D	328	HIS
1	D	415	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	SUG	A	1001	-	9,18,18	0.35	0	8,22,22	0.66	0
3	SUG	B	1002	-	9,18,18	0.37	0	8,22,22	0.24	0
3	SUG	C	1003	-	9,18,18	0.24	0	8,22,22	0.55	0
3	SUG	D	1004	-	9,18,18	0.48	0	8,22,22	0.42	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SUG	A	1001	-	-	0/12/20/20	0/0/0/0
3	SUG	B	1002	-	-	0/12/20/20	0/0/0/0
3	SUG	C	1003	-	-	0/12/20/20	0/0/0/0
3	SUG	D	1004	-	-	0/12/20/20	0/0/0/0

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.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1001	SUG	1	0
3	B	1002	SUG	4	0
3	D	1004	SUG	2	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	440/458 (96%)	-0.36	2 (0%) 91 91	10, 21, 36, 45	0
1	B	440/458 (96%)	-0.42	2 (0%) 91 91	10, 20, 35, 45	0
1	C	440/458 (96%)	-0.24	3 (0%) 89 88	11, 24, 37, 45	0
1	D	440/458 (96%)	-0.41	0 100 100	10, 22, 35, 45	0
All	All	1760/1832 (96%)	-0.36	7 (0%) 93 93	10, 22, 36, 45	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	95	HIS	3.2
1	A	4	TRP	3.1
1	B	313	ASP	2.6
1	A	313	ASP	2.4
1	C	2	ASN	2.4
1	C	3	ALA	2.3
1	B	91	ARG	2.2

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	SUG	C	1003	19/19	0.94	0.12	0.97	16,19,23,23	0
3	SUG	D	1004	19/19	0.97	0.09	-0.14	10,13,16,17	0
3	SUG	B	1002	19/19	0.97	0.10	-0.16	5,11,14,15	0
3	SUG	A	1001	19/19	0.97	0.09	-0.95	12,14,16,17	0
2	K	D	2004	1/1	0.91	0.10	-1.14	66,66,66,66	0
2	K	C	2003	1/1	0.79	0.10	-1.42	63,63,63,63	0
2	K	B	2002	1/1	0.97	0.08	-2.82	36,36,36,36	0
2	K	A	2001	1/1	0.99	0.04	-2.94	37,37,37,37	0

6.5 Other polymers

There are no such residues in this entry.