



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:29 PM GMT

PDB ID : 1KH1
Title : Crystal Structure of Thermus thermophilus HB8 Argininosuccinate Synthetase
Authors : goto, m.; Nakajima, Y; Hirotsu, k.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2001-11-29
Resolution : 2.30 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

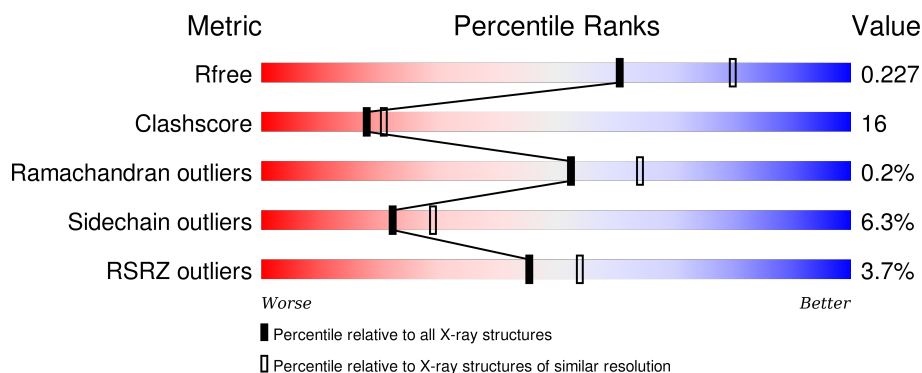
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	400	<div> <div>2%</div> <div>66%</div> <div>26%</div> <div>5%</div> </div>
1	B	400	<div> <div>%</div> <div>70%</div> <div>23%</div> <div>5%</div> </div>
1	C	400	<div> <div>5%</div> <div>70%</div> <div>23%</div> <div>• •</div> </div>
1	D	400	<div> <div>6%</div> <div>65%</div> <div>28%</div> <div>• •</div> </div>

2 Entry composition [i](#)

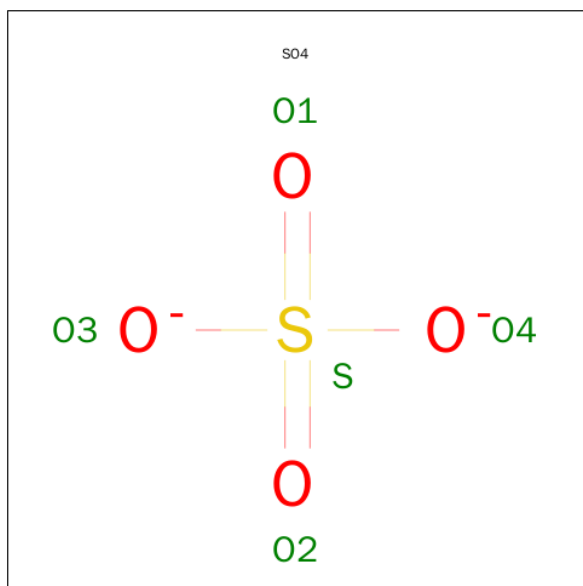
There are 3 unique types of molecules in this entry. The entry contains 12650 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Argininosuccinate Synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	380	Total	C	N	O	S	0	0	0
			3017	1931	526	551	9			
1	B	382	Total	C	N	O	S	0	0	0
			3034	1940	529	556	9			
1	C	386	Total	C	N	O	S	0	0	0
			3066	1963	533	561	9			
1	D	385	Total	C	N	O	S	0	0	0
			3055	1954	532	560	9			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	132	Total	O	0	0
			132	132		
3	B	124	Total	O	0	0
			124	124		
3	C	105	Total	O	0	0
			105	105		
3	D	97	Total	O	0	0
			97	97		

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.

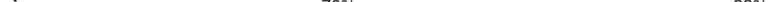
Chain A:

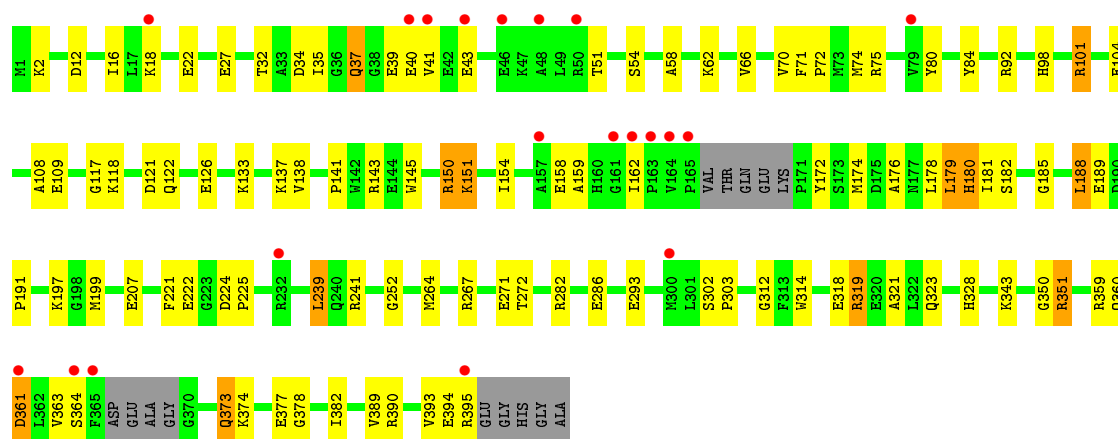
66% 26% 2% 5%

M1 K2 I3 V4 Y7 S8 G9 L11 I15 K21 Y24 A26 E27 V41 E42 E43 E46 R50 S54 A58 V66 D67 D68 F69 V70 M73 M74 R75 Y80 E81 Y84 L85 L86 I90 P93 K97 H98 V100 R101 E104 A109 A110 I111 G117 K118 G119 N120 D121 Q122 V123 R124 F125 G126 L127 T128 K133 I136 K137 V138 I139 R143 R144 H145 S146 F147 R150 K151 I154 A155 A156 A157 G161 H162 P163 P165 VAL TTR GLN GLU LYS P171 Y172 S173 M174 L179 H180 I181 S182 G185 L188 E189 P190 P191 E194 K197 V202 E214 F221 E222 A236 G239 H249 G252 V257 F261 R267 E271 T272 R282 E286 HIS E293 R298 S302 P303 Y315 A316 F317 E318 R319 E320 A321 L322 Q323 H328 R331 L339 N345 V346 Y347 V348 V349 G350 R351 Y358 R359 GLN ASP LEU VAL SER PHE ASP GLU GLY GLY G370 Y371 D372 Q373 Q383 R386 R390 V393 E394 R395 GLU GLY HIS GLY ALA

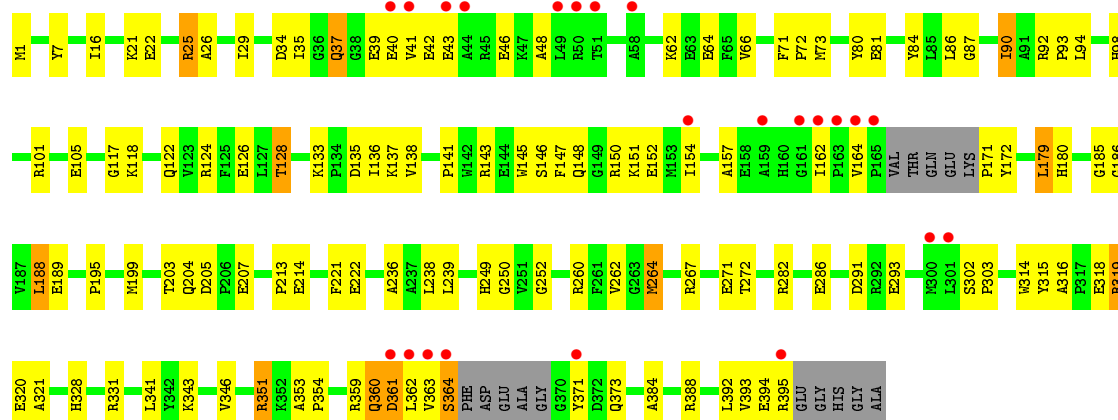
Chain B:

Amino Acid	Frequency (%)
M1	0.5
K2	0.5
A26	0.5
T32	0.5
A33	0.5
D34	0.5
I35	0.5
G36	0.5
Q37	0.5
E40	0.5
V41	0.5
E42	0.5
E46	0.5
A58	0.5
K62	0.5
S173	0.5
D68	0.5
P72	0.5
Y80	0.5
Y84	0.5
L85	0.5
L86	0.5
I90	0.5
A91	0.5
R92	0.5
P93	0.5
L94	0.5
K97	0.5
H98	0.5
I99	0.5
V100	0.5
R101	0.5
E104	0.5
E105	0.5
A108	0.5
E109	0.5
A110	0.5
K118	0.5
D121	0.5
Q122	0.5
V123	0.5
R124	0.5
F125	0.5
I126	0.5
I127	0.5
T128	0.5
I136	0.5
K137	0.5
V138	0.5
R143	0.5
E144	0.5
W145	0.5
K151	0.5
P165	0.5
VAL	0.5
THR	0.5
GLN	0.5
GLU	0.5
LYS	0.5
P171	0.5
Y172	0.5
S173	0.5
M174	0.5
L179	0.5
H180	0.5
I181	0.5
Y183	0.5
E184	0.5
G185	0.5
G186	0.5
V187	0.5
L188	0.5
P191	0.5
E194	0.5
P195	0.5
P196	0.5
M199	0.5
E214	0.5
Y215	0.5
V216	0.5
F221	0.5
E222	0.5
L238	0.5
L239	0.5
E249	0.5
F261	0.5
R267	0.5
Y270	0.5
E271	0.5
T272	0.5
R282	0.5
E286	0.5
S287	0.5
L288	0.5
D291	0.5
R292	0.5
E293	0.5
R298	0.5
D299	0.5
P300	0.5
K314	0.5
E318	0.5
R319	0.5
E320	0.5
A321	0.5
L322	0.5
Q323	0.5
E328	0.5
R338	0.5
V349	0.5
G350	0.5
R351	0.5
K352	0.5
A353	0.5
P354	0.5
K355	0.5
Y358	0.5
R359	0.5
Q360	0.5
D361	0.5
LEU	0.5
VAL	0.5
SER	0.5
PHE	0.5
ASP	0.5
GLU	0.5
ALA	0.5
GLY	0.5
G370	0.5
Q373	0.5
A376	0.5
F379	0.5
I380	0.5
R390	0.5
A391	0.5

Chain C:  5% 70% 23% . .



• Molecule 1: Argininosuccinate Synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	228.41Å 228.41Å 161.53Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.30 48.70 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.30) 89.2 (48.70-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 2.29Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.199 , 0.229 0.199 , 0.227	Depositor DCC
R_{free} test set	12258 reflections (9.98%)	DCC
Wilson B-factor (Å ²)	32.5	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 45.0	EDS
Estimated twinning fraction	0.019 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 124435 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12650	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.66% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: SO4

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	0.39	0/3085	0.66	4/4169 (0.1%)
1	B	0.39	0/3102	0.67	2/4192 (0.0%)
1	C	0.40	0/3135	0.66	3/4237 (0.1%)
1	D	0.39	0/3123	0.67	3/4221 (0.1%)
All	All	0.39	0/12445	0.66	12/16819 (0.1%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	221	PHE	CB-CA-C	-12.49	85.42	110.40
1	C	221	PHE	CB-CA-C	-11.59	87.22	110.40
1	A	221	PHE	CB-CA-C	-10.84	88.72	110.40
1	D	221	PHE	CB-CA-C	-10.63	89.13	110.40
1	C	222	GLU	N-CA-C	7.83	132.15	111.00
1	A	222	GLU	N-CA-C	7.54	131.35	111.00
1	D	222	GLU	N-CA-C	7.32	130.76	111.00
1	A	221	PHE	N-CA-C	-6.07	94.60	111.00
1	B	222	GLU	N-CA-C	5.94	127.03	111.00
1	C	221	PHE	N-CA-C	-5.72	95.54	111.00
1	D	221	PHE	N-CA-C	-5.68	95.66	111.00
1	A	257	VAL	N-CA-C	-5.13	97.15	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3017	0	3006	114	0
1	B	3034	0	3018	103	0
1	C	3066	0	3052	104	0
1	D	3055	0	3043	108	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	132	0	0	9	0
3	B	124	0	0	3	0
3	C	105	0	0	6	0
3	D	97	0	0	2	0
All	All	12650	0	12119	384	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 (384) 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:151:LYS:HD2	1:B:151:LYS:H	1.08	1.08
1:D:151:LYS:H	1:D:151:LYS:HD2	1.11	1.07
1:C:37:GLN:H	1:C:37:GLN:NE2	1.53	1.07
1:C:151:LYS:HD2	1:C:151:LYS:H	1.18	1.04
1:D:35:ILE:H	1:D:37:GLN:NE2	1.59	1.01
1:C:37:GLN:N	1:C:37:GLN:HE21	1.60	0.99
1:A:349:VAL:HG13	1:B:194:GLU:HB2	1.47	0.96
1:D:35:ILE:H	1:D:37:GLN:HE22	1.07	0.96
1:C:395:ARG:HG3	1:C:395:ARG:HH11	1.31	0.92
1:B:35:ILE:H	1:B:37:GLN:NE2	1.67	0.90
1:C:319:ARG:HH21	1:C:323:GLN:HE21	1.20	0.89
1:C:151:LYS:CD	1:C:151:LYS:H	1.83	0.89
1:D:393:VAL:C	1:D:395:ARG:H	1.75	0.88
1:A:150:ARG:HB3	1:A:151:LYS:HE3	1.55	0.88
1:C:37:GLN:H	1:C:37:GLN:HE21	0.92	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:151:LYS:HD2	1:C:151:LYS:N	1.91	0.86
1:B:151:LYS:CD	1:B:151:LYS:H	1.84	0.86
1:A:123:VAL:HG22	1:B:380:ILE:HD11	1.57	0.85
1:A:390:ARG:O	1:A:394:GLU:HG2	1.79	0.83
1:B:34:ASP:HB2	1:B:41:VAL:HG21	1.61	0.82
1:D:151:LYS:N	1:D:151:LYS:HD2	1.95	0.82
1:C:122:GLN:NE2	1:C:143:ARG:HH11	1.78	0.81
1:B:37:GLN:NE2	1:B:37:GLN:H	1.79	0.80
1:D:351:ARG:HD2	1:D:351:ARG:N	1.97	0.79
1:A:151:LYS:H	1:A:151:LYS:HE3	1.48	0.79
1:D:151:LYS:H	1:D:151:LYS:CD	1.90	0.78
1:B:1:MET:O	1:B:26:ALA:HB1	1.84	0.78
1:B:126:GLU:HG2	1:B:138:VAL:HG11	1.66	0.78
1:D:361:ASP:HA	1:D:364:SER:HB2	1.64	0.78
1:A:293:GLU:HB3	1:C:318:GLU:HG2	1.67	0.76
1:A:41:VAL:HG12	1:A:58:ALA:HB1	1.67	0.76
1:D:42:GLU:O	1:D:46:GLU:HG3	1.86	0.76
1:A:318:GLU:HG2	1:C:293:GLU:CB	2.16	0.76
1:A:293:GLU:CB	1:C:318:GLU:HG2	2.15	0.76
1:B:151:LYS:HD2	1:B:151:LYS:N	1.94	0.75
1:C:351:ARG:N	1:C:351:ARG:HD2	2.01	0.74
1:A:318:GLU:HG2	1:C:293:GLU:HB3	1.68	0.74
1:A:373:GLN:H	1:A:373:GLN:NE2	1.86	0.74
1:A:70:VAL:O	1:A:74:MET:HG3	1.86	0.74
1:C:122:GLN:HE21	1:C:143:ARG:HH11	1.34	0.73
1:D:171:PRO:O	1:D:186:GLY:HA3	1.87	0.73
1:B:351:ARG:N	1:B:351:ARG:HD2	2.02	0.73
1:B:359:ARG:HB3	1:B:361:ASP:OD1	1.88	0.73
1:D:393:VAL:C	1:D:395:ARG:N	2.40	0.73
1:A:98:HIS:HD2	1:A:101:ARG:NH1	1.85	0.73
1:C:319:ARG:HH21	1:C:323:GLN:NE2	1.87	0.72
1:A:351:ARG:N	1:A:351:ARG:HD2	2.04	0.72
1:A:109:GLU:O	1:A:137:LYS:HE2	1.88	0.72
1:D:35:ILE:N	1:D:37:GLN:HE22	1.85	0.71
1:B:100:VAL:O	1:B:104:GLU:HG3	1.90	0.71
1:D:126:GLU:HG2	1:D:138:VAL:HG11	1.73	0.70
1:B:293:GLU:HB3	1:D:318:GLU:HG2	1.71	0.70
1:D:395:ARG:NH1	1:D:395:ARG:HB3	2.06	0.70
1:B:293:GLU:CB	1:D:318:GLU:HG2	2.21	0.70
1:C:18:LYS:HG3	1:C:162:ILE:HD11	1.72	0.70
1:C:151:LYS:HE3	3:C:1356:HOH:O	1.90	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:395:ARG:CG	1:C:395:ARG:HH11	2.04	0.70
1:C:18:LYS:CG	1:C:162:ILE:HD11	2.21	0.70
1:A:123:VAL:HG22	1:B:380:ILE:CD1	2.20	0.69
1:D:179:LEU:O	1:D:272:THR:HG23	1.93	0.69
1:C:74:MET:HE2	3:C:1315:HOH:O	1.93	0.68
1:B:122:GLN:HE21	1:B:143:ARG:HH11	1.41	0.68
1:A:67:ARG:NH1	1:A:68:ASP:OD1	2.27	0.67
1:B:35:ILE:H	1:B:37:GLN:HE22	1.41	0.67
1:A:27:GLU:HG3	1:A:54:SER:CB	2.24	0.67
1:C:264:MET:HB2	1:D:362:LEU:HD13	1.76	0.67
1:B:41:VAL:HG12	1:B:58:ALA:HB1	1.77	0.67
1:B:314:TRP:O	1:B:319:ARG:HD3	1.94	0.67
1:D:282:ARG:O	1:D:286:GLU:HG3	1.95	0.67
1:A:93:PRO:HA	1:A:128:THR:HG21	1.77	0.66
1:D:40:GLU:HG3	1:D:43:GLU:HG3	1.76	0.66
1:D:40:GLU:CG	1:D:43:GLU:HG3	2.26	0.66
1:D:395:ARG:HB3	1:D:395:ARG:CZ	2.26	0.66
1:D:98:HIS:HD2	1:D:101:ARG:HH12	1.44	0.66
1:B:32:THR:HG22	1:B:41:VAL:HG13	1.76	0.66
1:B:318:GLU:HG2	1:D:293:GLU:HB3	1.78	0.66
1:C:32:THR:CG2	1:C:41:VAL:HG13	2.27	0.65
1:D:124:ARG:O	1:D:128:THR:HG23	1.96	0.65
1:A:151:LYS:CE	1:A:151:LYS:H	2.08	0.65
1:A:293:GLU:HB2	1:C:318:GLU:CG	2.25	0.65
1:D:98:HIS:HD2	1:D:101:ARG:NH1	1.95	0.65
1:A:349:VAL:CG1	1:B:194:GLU:HB2	2.25	0.64
1:B:37:GLN:HE21	1:B:37:GLN:H	1.44	0.64
1:B:393:VAL:C	1:B:395:ARG:H	2.01	0.64
1:A:133:LYS:HG2	1:A:136:ILE:HB	1.79	0.63
1:B:35:ILE:N	1:B:37:GLN:NE2	2.42	0.63
1:D:98:HIS:CD2	1:D:101:ARG:NH1	2.66	0.63
1:A:151:LYS:CD	1:A:151:LYS:H	2.12	0.63
1:A:21:LYS:HD3	3:A:1313:HOH:O	1.99	0.63
1:C:118:LYS:HG2	3:D:1428:HOH:O	1.98	0.63
1:D:117:GLY:HA2	1:D:122:GLN:HE22	1.63	0.63
1:D:133:LYS:HD3	1:D:136:ILE:HB	1.80	0.63
1:C:41:VAL:HG12	1:C:58:ALA:HB1	1.79	0.62
1:C:35:ILE:H	1:C:37:GLN:NE2	1.98	0.62
1:D:1:MET:O	1:D:26:ALA:HB1	2.00	0.61
1:A:126:GLU:HG2	1:A:138:VAL:HG11	1.81	0.61
1:C:302:SER:HB2	1:C:303:PRO:HD3	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:189:GLU:O	1:C:191:PRO:HD3	2.00	0.60
1:B:293:GLU:HB2	1:D:318:GLU:CG	2.31	0.60
1:A:41:VAL:CG1	1:A:58:ALA:HB1	2.30	0.60
1:B:1:MET:HB3	1:B:26:ALA:HB2	1.82	0.60
1:A:214:GLU:OE1	1:A:249:HIS:HE1	1.84	0.60
1:D:360:GLN:HA	1:D:363:VAL:HG22	1.84	0.60
1:D:90:ILE:HD11	1:D:179:LEU:HD12	1.83	0.60
1:C:16:ILE:HG23	1:C:141:PRO:HG2	1.84	0.59
1:D:117:GLY:HA2	1:D:122:GLN:NE2	2.17	0.59
1:B:97:LYS:NZ	3:B:1116:HOH:O	2.35	0.59
1:B:359:ARG:O	1:B:361:ASP:N	2.35	0.59
1:A:331:ARG:HH11	1:A:331:ARG:CB	2.16	0.59
1:A:318:GLU:CG	1:C:293:GLU:HB2	2.32	0.59
1:D:37:GLN:NE2	1:D:37:GLN:H	2.00	0.58
1:B:314:TRP:CE3	1:B:319:ARG:HD2	2.38	0.58
1:A:370:GLY:N	3:A:1105:HOH:O	2.35	0.58
1:A:249:HIS:HD2	3:A:1155:HOH:O	1.85	0.58
1:B:32:THR:CG2	1:B:41:VAL:HG13	2.33	0.58
1:B:376:ALA:O	1:B:380:ILE:HG12	2.02	0.58
1:D:124:ARG:O	1:D:128:THR:CG2	2.51	0.58
1:C:12:ASP:OD1	1:C:150:ARG:NH2	2.37	0.58
1:D:314:TRP:O	1:D:319:ARG:HD3	2.04	0.57
1:C:185:GLY:O	1:C:188:LEU:HB2	2.04	0.57
1:B:35:ILE:N	1:B:37:GLN:HE22	2.00	0.57
1:A:282:ARG:HD2	3:A:1115:HOH:O	2.04	0.57
1:A:395:ARG:NH1	3:A:1301:HOH:O	2.37	0.57
1:A:349:VAL:HG12	3:A:1418:HOH:O	2.05	0.57
1:D:395:ARG:NH1	1:D:395:ARG:CB	2.68	0.57
1:A:393:VAL:C	1:A:395:ARG:H	2.07	0.57
1:B:318:GLU:HG2	1:D:293:GLU:CB	2.35	0.57
1:D:122:GLN:HE21	1:D:143:ARG:HH11	1.53	0.57
1:D:35:ILE:N	1:D:37:GLN:NE2	2.43	0.56
1:B:171:PRO:O	1:B:186:GLY:HA3	2.04	0.56
1:C:122:GLN:HE22	1:C:143:ARG:HD3	1.70	0.56
1:C:395:ARG:NH1	1:C:395:ARG:HG3	2.11	0.56
1:B:214:GLU:OE1	1:B:249:HIS:HE1	1.89	0.56
1:C:32:THR:HG22	1:C:41:VAL:HG13	1.87	0.56
1:C:282:ARG:O	1:C:286:GLU:HG3	2.06	0.56
1:A:90:ILE:HD11	1:A:179:LEU:HD12	1.86	0.56
1:B:282:ARG:O	1:B:286:GLU:HG3	2.05	0.56
1:B:42:GLU:O	1:B:46:GLU:HG3	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:GLU:HG2	1:C:293:GLU:HB2	1.84	0.56
1:D:162:ILE:O	1:D:164:VAL:HG13	2.06	0.56
1:B:122:GLN:NE2	1:B:143:ARG:HH11	2.02	0.56
1:B:93:PRO:HA	1:B:128:THR:HG21	1.88	0.56
1:D:21:LYS:O	1:D:25:ARG:HA	2.05	0.56
1:D:316:ALA:HB1	1:D:318:GLU:OE2	2.05	0.55
1:C:150:ARG:O	1:C:154:ILE:HG13	2.06	0.55
1:A:302:SER:HB2	1:A:303:PRO:HD3	1.87	0.55
1:C:319:ARG:NH2	1:C:323:GLN:HE21	1.96	0.55
1:B:318:GLU:CG	1:D:293:GLU:HB2	2.37	0.55
1:D:148:GLN:HB2	1:D:152:GLU:OE2	2.06	0.55
1:A:358:TYR:CE2	1:B:191:PRO:HD3	2.40	0.55
1:C:122:GLN:HE21	1:C:143:ARG:NH1	2.02	0.55
1:A:293:GLU:HB2	1:C:318:GLU:HG2	1.84	0.55
1:D:80:TYR:CD2	1:D:81:GLU:HG3	2.42	0.55
1:C:179:LEU:O	1:C:272:THR:HG23	2.07	0.54
1:A:98:HIS:CD2	1:A:101:ARG:NH1	2.73	0.54
1:D:214:GLU:OE1	1:D:249:HIS:HE1	1.90	0.54
1:A:372:ASP:HB2	1:A:373:GLN:HE21	1.70	0.54
1:A:66:VAL:HG21	1:A:236:ALA:HA	1.90	0.54
1:C:343:LYS:HD2	1:D:343:LYS:CD	2.37	0.54
1:B:118:LYS:HD2	1:B:118:LYS:H	1.72	0.54
1:D:185:GLY:O	1:D:188:LEU:HB2	2.07	0.54
1:A:331:ARG:HB3	1:A:331:ARG:NH1	2.22	0.54
1:A:86:LEU:O	1:A:90:ILE:HG13	2.07	0.54
1:A:319:ARG:HH21	1:A:323:GLN:NE2	2.06	0.54
1:A:27:GLU:HG3	1:A:54:SER:HB3	1.88	0.54
1:C:70:VAL:O	1:C:74:MET:HG3	2.08	0.53
1:B:249:HIS:HD2	3:B:1427:HOH:O	1.89	0.53
1:C:343:LYS:HD2	1:D:343:LYS:HD3	1.90	0.53
1:C:364:SER:O	1:D:118:LYS:NZ	2.35	0.53
1:D:262:VAL:CG1	1:D:262:VAL:O	2.55	0.53
1:B:379:PHE:HB2	1:C:382:ILE:HG13	1.90	0.53
1:A:120:ASN:HD21	1:A:261:PHE:H	1.57	0.53
1:D:392:LEU:O	1:D:395:ARG:HA	2.08	0.53
1:B:393:VAL:C	1:B:395:ARG:N	2.61	0.53
1:B:2:LYS:HB2	1:B:108:ALA:HA	1.90	0.53
1:B:34:ASP:HA	1:B:37:GLN:HE22	1.73	0.53
1:C:176:ALA:CB	1:C:181:ILE:HG12	2.39	0.52
1:B:37:GLN:HE21	1:B:37:GLN:N	2.07	0.52
1:C:373:GLN:NE2	1:C:373:GLN:H	2.07	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:390:ARG:O	1:B:394:GLU:HG2	2.10	0.52
1:C:394:GLU:O	1:C:395:ARG:HB2	2.10	0.52
1:B:124:ARG:O	1:B:128:THR:HG23	2.10	0.52
1:B:98:HIS:HD2	1:B:101:ARG:NH1	2.07	0.52
1:C:34:ASP:HB2	1:C:41:VAL:HG21	1.90	0.52
1:C:104:GLU:OE2	1:C:133:LYS:HD3	2.10	0.52
1:A:137:LYS:N	1:A:137:LYS:HD3	2.24	0.52
1:D:150:ARG:HG3	1:D:154:ILE:HD11	1.92	0.52
1:B:80:TYR:HB3	1:B:84:TYR:HB3	1.91	0.52
1:D:87:GLY:HA2	1:D:179:LEU:HD13	1.91	0.51
1:D:148:GLN:H	1:D:152:GLU:CD	2.12	0.51
1:C:37:GLN:HB2	1:C:39:GLU:HG2	1.92	0.51
1:B:319:ARG:HH21	1:B:323:GLN:HE21	1.56	0.51
1:C:2:LYS:HB2	1:C:108:ALA:HA	1.92	0.51
1:B:86:LEU:O	1:B:90:ILE:HG13	2.11	0.51
1:D:393:VAL:O	1:D:395:ARG:N	2.44	0.51
1:C:109:GLU:O	1:C:137:LYS:HE2	2.11	0.51
1:C:18:LYS:HG2	1:C:162:ILE:HD11	1.92	0.51
1:A:40:GLU:HG3	1:A:43:GLU:HG3	1.93	0.51
1:D:93:PRO:HA	1:D:128:THR:HG21	1.93	0.51
1:A:21:LYS:HD2	1:A:26:ALA:O	2.11	0.50
1:A:97:LYS:HD2	1:A:315:TYR:OH	2.11	0.50
1:C:373:GLN:HG2	1:C:374:LYS:N	2.27	0.50
1:B:179:LEU:HD22	1:B:272:THR:HG23	1.93	0.50
1:A:293:GLU:CB	1:C:318:GLU:CG	2.86	0.50
1:B:123:VAL:HG12	1:B:127:LEU:HD12	1.94	0.50
1:D:360:GLN:O	1:D:364:SER:HB2	2.11	0.50
1:A:373:GLN:H	1:A:373:GLN:HE21	1.57	0.50
1:D:260:ARG:HD2	1:D:264:MET:SD	2.52	0.50
1:B:359:ARG:C	1:B:361:ASP:H	2.15	0.50
1:D:84:TYR:CE2	1:D:86:LEU:HA	2.46	0.50
1:D:252:GLY:O	1:D:271:GLU:HA	2.11	0.50
1:C:378:GLY:O	1:C:382:ILE:HG12	2.11	0.49
1:D:351:ARG:HD2	1:D:351:ARG:H	1.74	0.49
1:C:390:ARG:O	1:C:394:GLU:HG2	2.13	0.49
1:A:319:ARG:HH21	1:A:323:GLN:HE21	1.60	0.49
1:C:27:GLU:HG3	1:C:54:SER:OG	2.12	0.49
1:B:98:HIS:CD2	1:B:101:ARG:NH1	2.80	0.49
1:C:176:ALA:HB2	1:C:181:ILE:HG12	1.94	0.49
1:C:241:ARG:NH1	3:C:1267:HOH:O	2.43	0.49
1:C:51:THR:HG22	1:C:51:THR:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:393:VAL:C	1:A:395:ARG:N	2.66	0.49
1:B:338:ARG:HD3	3:B:1416:HOH:O	2.12	0.49
1:B:122:GLN:HE22	1:B:143:ARG:HD3	1.78	0.49
1:B:318:GLU:CG	1:D:293:GLU:CB	2.91	0.49
1:A:24:TYR:OH	1:A:144:GLU:OE2	2.28	0.48
1:D:66:VAL:HG21	1:D:236:ALA:HA	1.95	0.48
1:D:101:ARG:HG2	1:D:105:GLU:OE2	2.14	0.48
1:D:328:HIS:CE1	1:D:331:ARG:HH22	2.31	0.48
1:D:384:ALA:O	1:D:388:ARG:HG3	2.13	0.48
1:B:36:GLY:N	1:B:37:GLN:HE21	2.12	0.48
1:A:84:TYR:CE2	1:A:86:LEU:HA	2.49	0.48
1:A:339:LEU:N	1:A:339:LEU:HD12	2.29	0.48
1:C:178:LEU:HB3	1:C:239:LEU:HD13	1.96	0.48
1:C:180:HIS:ND1	1:C:180:HIS:C	2.67	0.48
1:D:395:ARG:CB	1:D:395:ARG:HH11	2.26	0.48
1:D:16:ILE:HG23	1:D:141:PRO:HG2	1.95	0.48
1:A:146:SER:O	1:A:147:PHE:C	2.53	0.47
1:C:174:MET:HA	1:C:182:SER:O	2.14	0.47
1:B:180:HIS:ND1	1:B:180:HIS:C	2.68	0.47
1:D:314:TRP:CE3	1:D:319:ARG:HD2	2.49	0.47
1:A:179:LEU:HD22	1:A:272:THR:HG23	1.97	0.47
1:A:98:HIS:CD2	1:A:101:ARG:HH11	2.33	0.47
1:A:120:ASN:ND2	1:A:261:PHE:H	2.12	0.47
1:C:395:ARG:CG	1:C:395:ARG:NH1	2.68	0.47
1:B:179:LEU:O	1:B:272:THR:HG23	2.14	0.47
1:C:180:HIS:HB2	1:C:271:GLU:O	2.14	0.47
1:A:316:ALA:HB1	1:A:318:GLU:OE2	2.15	0.47
1:C:252:GLY:O	1:C:271:GLU:HA	2.14	0.47
1:D:80:TYR:CE2	1:D:81:GLU:HG3	2.49	0.47
1:B:185:GLY:O	1:B:188:LEU:HB2	2.15	0.47
1:D:22:GLU:HA	1:D:25:ARG:NH1	2.30	0.47
1:A:3:ILE:HD13	1:A:139:ILE:HD12	1.97	0.47
1:C:389:VAL:O	1:C:393:VAL:HG23	2.15	0.46
1:C:75:ARG:NH2	3:C:1156:HOH:O	2.48	0.46
1:C:101:ARG:NH2	3:C:1140:HOH:O	2.34	0.46
1:A:154:ILE:HG22	1:A:155:ALA:N	2.31	0.46
1:A:121:ASP:HA	1:A:124:ARG:HD2	1.96	0.46
1:B:355:LYS:HE3	1:C:393:VAL:HG12	1.96	0.46
1:A:21:LYS:O	1:A:25:ARG:HA	2.16	0.46
1:B:359:ARG:C	1:B:361:ASP:N	2.68	0.46
1:C:314:TRP:O	1:C:319:ARG:HD3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:84:TYR:CZ	1:D:86:LEU:HA	2.50	0.46
1:B:328:HIS:CD2	1:D:321:ALA:HA	2.50	0.46
1:D:146:SER:O	1:D:147:PHE:C	2.53	0.46
1:C:35:ILE:H	1:C:37:GLN:HE22	1.64	0.46
1:C:98:HIS:HD2	1:C:101:ARG:NH1	2.13	0.46
1:B:318:GLU:HG3	1:D:293:GLU:HB2	1.97	0.46
1:C:126:GLU:HG2	1:C:138:VAL:HG11	1.97	0.46
1:D:203:THR:HB	1:D:250:GLY:HA2	1.98	0.46
3:A:1265:HOH:O	1:B:191:PRO:HG2	2.16	0.45
1:A:189:GLU:OE2	1:B:360:GLN:OE1	2.34	0.45
1:B:34:ASP:HB2	1:B:41:VAL:CG2	2.39	0.45
1:B:293:GLU:HB2	1:D:318:GLU:HG2	1.92	0.45
1:B:182:SER:HB2	1:B:270:TYR:CE2	2.52	0.45
1:C:35:ILE:N	1:C:37:GLN:HE22	2.15	0.45
1:D:29:ILE:N	1:D:29:ILE:HD12	2.31	0.45
1:C:92:ARG:HE	1:C:121:ASP:CG	2.19	0.45
1:C:374:LYS:HE2	1:C:377:GLU:OE2	2.17	0.45
1:A:373:GLN:H	1:A:373:GLN:CD	2.16	0.45
1:A:11:LEU:O	1:A:15:ILE:HG13	2.17	0.45
1:D:150:ARG:HG3	1:D:154:ILE:CD1	2.46	0.45
1:B:110:ALA:HB2	1:B:137:LYS:HB3	1.98	0.45
1:A:151:LYS:N	1:A:151:LYS:HE3	2.25	0.44
1:A:331:ARG:NH1	1:A:331:ARG:CB	2.79	0.44
1:A:80:TYR:HB3	1:A:84:TYR:HB3	1.98	0.44
1:A:11:LEU:HD21	1:A:154:ILE:HD12	1.98	0.44
1:D:302:SER:HB2	1:D:303:PRO:HD3	1.99	0.44
1:C:34:ASP:HA	1:C:37:GLN:HE22	1.82	0.44
1:D:128:THR:HB	1:D:315:TYR:HD1	1.83	0.44
1:B:174:MET:HG2	1:B:183:TYR:CD2	2.53	0.44
1:A:191:PRO:HD3	1:B:358:TYR:CE2	2.52	0.44
1:A:100:VAL:O	1:A:104:GLU:HG3	2.16	0.44
1:D:204:GLN:HG2	1:D:205:ASP:N	2.32	0.44
1:A:194:GLU:HB2	1:B:349:VAL:HB	1.99	0.44
1:B:126:GLU:CG	1:B:138:VAL:HG11	2.42	0.44
1:A:50:ARG:NH1	3:A:1289:HOH:O	2.50	0.44
1:A:252:GLY:O	1:A:271:GLU:HA	2.18	0.44
1:A:328:HIS:CD2	1:C:321:ALA:HA	2.53	0.44
1:A:257:VAL:HG11	1:B:288:LEU:HD23	2.00	0.44
1:C:66:VAL:HA	1:C:70:VAL:HB	2.00	0.44
1:D:133:LYS:HE2	1:D:135:ASP:OD2	2.18	0.44
1:B:321:ALA:HA	1:D:328:HIS:CD2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:GLU:HG3	1:A:54:SER:OG	2.17	0.43
1:B:172:TYR:CZ	1:B:199:MET:HG3	2.53	0.43
1:C:360:GLN:OE1	1:D:189:GLU:OE2	2.36	0.43
1:B:92:ARG:HB2	1:B:93:PRO:HD3	2.00	0.43
1:B:123:VAL:HG12	1:B:127:LEU:CD1	2.49	0.43
1:A:197:LYS:HG2	1:A:197:LYS:H	1.57	0.43
1:C:117:GLY:HA3	1:D:373:GLN:OE1	2.18	0.43
1:D:7:TYR:CD2	1:D:48:ALA:HB2	2.53	0.43
1:C:312:GLY:HA2	3:C:1083:HOH:O	2.19	0.43
1:D:316:ALA:O	1:D:320:GLU:HG3	2.19	0.43
1:D:64:GLU:OE1	1:D:98:HIS:HE1	2.01	0.43
1:D:154:ILE:O	1:D:157:ALA:HB3	2.19	0.43
1:D:341:LEU:HD22	1:D:346:VAL:HG22	2.00	0.43
1:B:122:GLN:NE2	1:B:143:ARG:HD3	2.33	0.43
1:C:109:GLU:HG2	1:C:137:LYS:CE	2.49	0.43
1:A:122:GLN:OE1	1:A:143:ARG:HD3	2.19	0.43
1:C:37:GLN:N	1:C:37:GLN:NE2	2.33	0.43
1:D:73:MET:CE	1:D:90:ILE:HG12	2.48	0.43
1:A:185:GLY:HA2	1:A:189:GLU:HG3	2.01	0.43
1:A:179:LEU:O	1:A:272:THR:HG23	2.18	0.43
1:A:4:VAL:O	1:A:111:ILE:HA	2.19	0.43
1:B:37:GLN:NE2	1:B:37:GLN:N	2.58	0.42
1:D:353:ALA:HA	1:D:354:PRO:HD3	1.90	0.42
1:D:92:ARG:HB2	1:D:93:PRO:HD3	2.01	0.42
1:A:90:ILE:CD1	1:A:179:LEU:HD12	2.49	0.42
1:B:101:ARG:O	1:B:105:GLU:HG3	2.19	0.42
1:A:174:MET:HA	1:A:182:SER:O	2.19	0.42
1:B:136:ILE:HG13	1:B:137:LYS:N	2.34	0.42
1:C:172:TYR:CZ	1:C:199:MET:HG3	2.54	0.42
1:A:321:ALA:HA	1:C:328:HIS:CD2	2.55	0.42
1:A:124:ARG:O	1:A:128:THR:HG23	2.19	0.42
1:A:331:ARG:HH11	1:A:331:ARG:HB2	1.84	0.42
1:C:360:GLN:O	1:C:363:VAL:HG22	2.20	0.42
1:B:238:LEU:HD23	1:B:238:LEU:O	2.19	0.42
1:A:15:ILE:HG23	1:A:156:TYR:CD2	2.55	0.42
1:D:205:ASP:OD2	1:D:207:GLU:HB2	2.20	0.42
1:A:7:TYR:CE2	1:A:9:GLY:HA2	2.54	0.42
1:A:85:LEU:HA	1:A:85:LEU:HD23	1.83	0.42
1:A:350:GLY:C	1:A:351:ARG:HD2	2.38	0.42
1:A:117:GLY:HA3	1:B:373:GLN:NE2	2.35	0.42
1:D:71:PHE:HB2	1:D:72:PRO:HD3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:350:GLY:C	1:C:351:ARG:HD2	2.40	0.42
1:D:171:PRO:HG2	1:D:172:TYR:H	1.84	0.41
1:B:238:LEU:HD23	1:B:238:LEU:C	2.40	0.41
1:D:34:ASP:HB2	1:D:41:VAL:HG11	2.02	0.41
1:B:351:ARG:N	1:B:351:ARG:CD	2.80	0.41
1:A:181:ILE:HD12	1:A:202:MET:HB2	2.02	0.41
1:C:80:TYR:HB3	1:C:84:TYR:HB3	2.02	0.41
1:D:359:ARG:NH2	1:D:361:ASP:OD1	2.52	0.41
1:C:343:LYS:HD2	1:D:343:LYS:HD2	2.02	0.41
1:C:224:ASP:HA	1:C:225:PRO:HD3	1.91	0.41
1:B:353:ALA:HA	1:B:354:PRO:HD3	1.91	0.41
1:D:360:GLN:O	1:D:364:SER:N	2.53	0.41
1:C:40:GLU:O	1:C:43:GLU:N	2.50	0.41
1:A:345:ASN:HB3	1:A:347:TYR:CE1	2.56	0.41
1:A:282:ARG:O	1:A:286:GLU:HG3	2.21	0.41
1:B:298:ARG:HD2	1:B:298:ARG:C	2.41	0.41
1:B:68:ASP:O	1:B:72:PRO:HG2	2.21	0.41
1:B:350:GLY:C	1:B:351:ARG:HD2	2.41	0.41
1:A:386:ARG:HA	1:C:264:MET:SD	2.61	0.41
1:B:392:LEU:O	1:B:395:ARG:HA	2.21	0.41
1:A:383:GLN:NE2	1:B:261:PHE:CZ	2.89	0.41
1:A:383:GLN:NE2	1:B:261:PHE:HZ	2.19	0.41
1:A:80:TYR:CD2	1:A:81:GLU:HG3	2.56	0.41
1:C:314:TRP:O	1:C:319:ARG:CD	2.69	0.40
1:A:73:MET:CE	1:A:90:ILE:HG12	2.51	0.40
1:A:154:ILE:HG13	1:A:164:VAL:HG11	2.02	0.40
1:A:185:GLY:O	1:A:188:LEU:HB2	2.21	0.40
1:C:359:ARG:NE	1:C:361:ASP:OD1	2.54	0.40
1:D:238:LEU:C	1:D:238:LEU:HD23	2.41	0.40
1:C:207:GLU:OE1	1:D:213:PRO:HG2	2.21	0.40
1:C:71:PHE:HB2	1:C:72:PRO:HD3	2.03	0.40
1:B:179:LEU:HD22	1:B:272:THR:CG2	2.50	0.40
1:A:180:HIS:HB2	1:A:271:GLU:O	2.22	0.40
1:A:298:ARG:C	1:A:298:ARG:HD2	2.41	0.40
1:A:118:LYS:H	1:A:118:LYS:HD2	1.87	0.40
1:D:195:PRO:HB3	1:D:199:MET:CE	2.51	0.40
1:D:359:ARG:NH2	3:D:1096:HOH:O	2.53	0.40
1:C:151:LYS:CD	1:C:151:LYS:N	2.58	0.40
1:A:154:ILE:CG2	1:A:155:ALA:N	2.84	0.40
1:B:196:PRO:HG2	1:B:199:MET:HB2	2.04	0.40
1:A:75:ARG:NH2	3:A:1346:HOH:O	2.54	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	374/400 (94%)	360 (96%)	13 (4%)	1 (0%)	46	57
1	B	376/400 (94%)	362 (96%)	13 (4%)	1 (0%)	46	57
1	C	380/400 (95%)	365 (96%)	14 (4%)	1 (0%)	46	57
1	D	379/400 (95%)	365 (96%)	14 (4%)	0	100	100
All	All	1509/1600 (94%)	1452 (96%)	54 (4%)	3 (0%)	52	64

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	360	GLN
1	A	147	PHE
1	C	159	ALA

5.3.2 Protein sidechains [i](#)

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	304/319 (95%)	289 (95%)	15 (5%)	31	41
1	B	306/319 (96%)	285 (93%)	21 (7%)	19	24
1	C	310/319 (97%)	292 (94%)	18 (6%)	25	33
1	D	309/319 (97%)	286 (93%)	23 (7%)	17	21

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1229/1276 (96%)	1152 (94%)	77 (6%)	22	29

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	ILE
1	A	101	ARG
1	A	118	LYS
1	A	128	THR
1	A	145	TRP
1	A	151	LYS
1	A	154	ILE
1	A	172	TYR
1	A	179	LEU
1	A	180	HIS
1	A	239	LEU
1	A	267	ARG
1	A	319	ARG
1	A	351	ARG
1	A	373	GLN
1	B	37	GLN
1	B	62	LYS
1	B	90	ILE
1	B	94	LEU
1	B	118	LYS
1	B	121	ASP
1	B	128	THR
1	B	145	TRP
1	B	151	LYS
1	B	165	PRO
1	B	179	LEU
1	B	180	HIS
1	B	184	GLU
1	B	216	VAL
1	B	239	LEU
1	B	267	ARG
1	B	291	ASP
1	B	319	ARG
1	B	351	ARG
1	B	355	LYS
1	B	361	ASP
1	C	22	GLU

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Mol	Chain	Res	Type
1	C	37	GLN
1	C	62	LYS
1	C	101	ARG
1	C	145	TRP
1	C	150	ARG
1	C	151	LYS
1	C	158	GLU
1	C	179	LEU
1	C	180	HIS
1	C	188	LEU
1	C	197	LYS
1	C	239	LEU
1	C	267	ARG
1	C	319	ARG
1	C	351	ARG
1	C	361	ASP
1	C	373	GLN
1	D	25	ARG
1	D	37	GLN
1	D	39	GLU
1	D	62	LYS
1	D	90	ILE
1	D	94	LEU
1	D	128	THR
1	D	137	LYS
1	D	145	TRP
1	D	179	LEU
1	D	180	HIS
1	D	188	LEU
1	D	239	LEU
1	D	264	MET
1	D	267	ARG
1	D	291	ASP
1	D	319	ARG
1	D	351	ARG
1	D	360	GLN
1	D	361	ASP
1	D	364	SER
1	D	371	TYR
1	D	394	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	HIS
1	A	249	HIS
1	A	323	GLN
1	A	373	GLN
1	B	37	GLN
1	B	98	HIS
1	B	122	GLN
1	B	249	HIS
1	B	323	GLN
1	B	360	GLN
1	C	37	GLN
1	C	98	HIS
1	C	122	GLN
1	C	323	GLN
1	C	360	GLN
1	C	373	GLN
1	D	37	GLN
1	D	98	HIS
1	D	122	GLN
1	D	148	GLN
1	D	249	HIS
1	D	297	GLN
1	D	323	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	510	-	4,4,4	0.25	0	6,6,6	0.15	0
2	SO4	B	1510	-	4,4,4	0.26	0	6,6,6	0.22	0
2	SO4	C	2510	-	4,4,4	0.42	0	6,6,6	0.07	0
2	SO4	D	3510	-	4,4,4	0.38	0	6,6,6	0.25	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	A	510	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1510	-	-	0/0/0/0	0/0/0/0
2	SO4	C	2510	-	-	0/0/0/0	0/0/0/0
2	SO4	D	3510	-	-	0/0/0/0	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.

No monomer is involved in short contacts.

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	380/400 (95%)	0.02	9 (2%) 62 71	20, 31, 54, 70	0
1	B	382/400 (95%)	-0.09	5 (1%) 79 84	22, 31, 48, 83	0
1	C	386/400 (96%)	0.17	20 (5%) 31 39	23, 34, 60, 82	0
1	D	385/400 (96%)	0.20	23 (5%) 25 33	20, 34, 59, 82	0
All	All	1533/1600 (95%)	0.08	57 (3%) 45 54	20, 32, 56, 83	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	365	PHE	4.7
1	C	165	PRO	4.4
1	D	165	PRO	4.2
1	B	360	GLN	4.1
1	D	163	PRO	4.0
1	D	164	VAL	3.9
1	D	40	GLU	3.3
1	D	362	LEU	3.3
1	D	361	ASP	3.2
1	D	371	TYR	3.2
1	A	154	ILE	3.1
1	A	157	ALA	3.1
1	B	361	ASP	3.1
1	D	41	VAL	3.1
1	C	157	ALA	2.9
1	B	165	PRO	2.9
1	C	50	ARG	2.9
1	A	162	ILE	2.8
1	D	154	ILE	2.8
1	C	43	GLU	2.8
1	D	161	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	163	PRO	2.8
1	D	300	MET	2.7
1	D	44	ALA	2.6
1	D	364	SER	2.6
1	A	41	VAL	2.6
1	A	165	PRO	2.6
1	C	361	ASP	2.5
1	C	161	GLY	2.5
1	D	159	ALA	2.5
1	C	41	VAL	2.5
1	D	363	VAL	2.5
1	C	40	GLU	2.5
1	A	395	ARG	2.4
1	C	395	ARG	2.4
1	D	50	ARG	2.4
1	C	162	ILE	2.4
1	C	364	SER	2.3
1	B	40	GLU	2.2
1	D	49	LEU	2.2
1	C	46	GLU	2.2
1	C	232	ARG	2.2
1	C	163	PRO	2.2
1	D	395	ARG	2.2
1	D	51	THR	2.2
1	D	58	ALA	2.2
1	B	300	MET	2.1
1	C	48	ALA	2.1
1	A	161	GLY	2.1
1	C	79	VAL	2.1
1	C	164	VAL	2.1
1	D	162	ILE	2.1
1	C	300	MET	2.1
1	D	43	GLU	2.1
1	C	18	LYS	2.1
1	A	46	GLU	2.0
1	D	301	LEU	2.0

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 [i](#)

There are no carbohydrates 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. 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(\AA^2)	Q<0.9
2	SO4	A	510	5/5	0.98	0.15	0.24	43,44,46,46	0
2	SO4	B	1510	5/5	0.99	0.12	0.20	41,44,47,49	0
2	SO4	D	3510	5/5	0.99	0.11	-1.06	45,49,51,54	0
2	SO4	C	2510	5/5	0.98	0.11	-1.44	50,51,53,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.