



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:39 AM UTC

PDB ID : 1KH1 / pdb_00001kh1
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

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

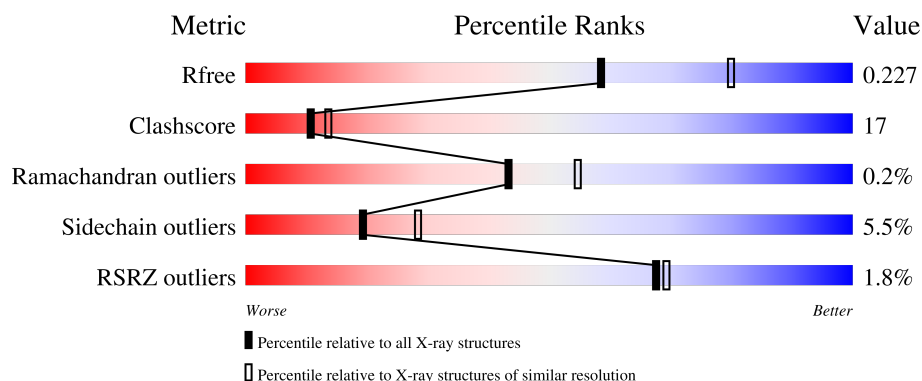
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	400	2% 65% 26% 5%
1	B	400	2% 68% 24%
1	C	400	2% 68% 24% 5%
1	D	400	2% 62% 30%

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 (CCD ID: 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		

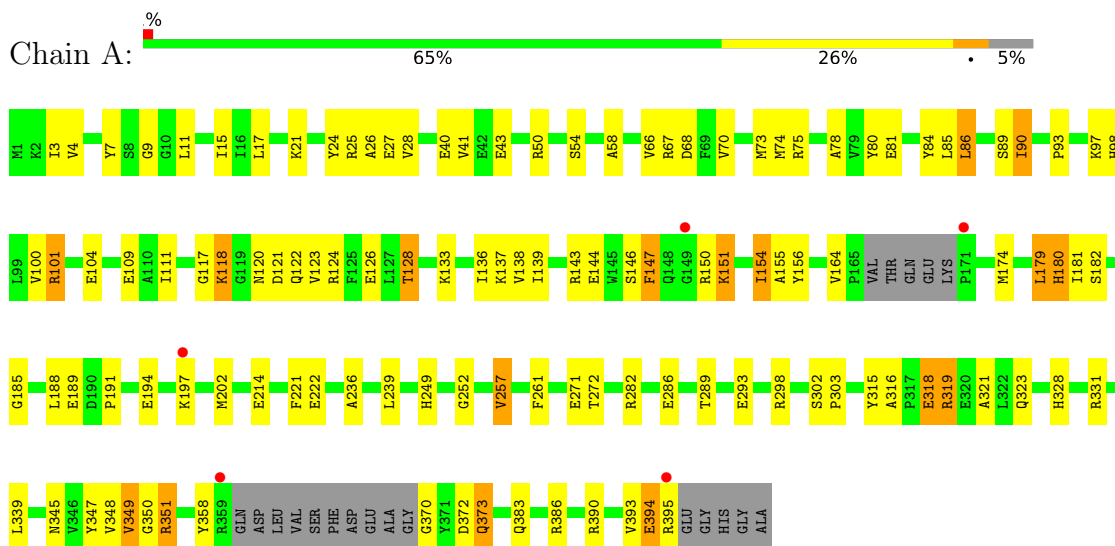
- Molecule 3 is water.

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		

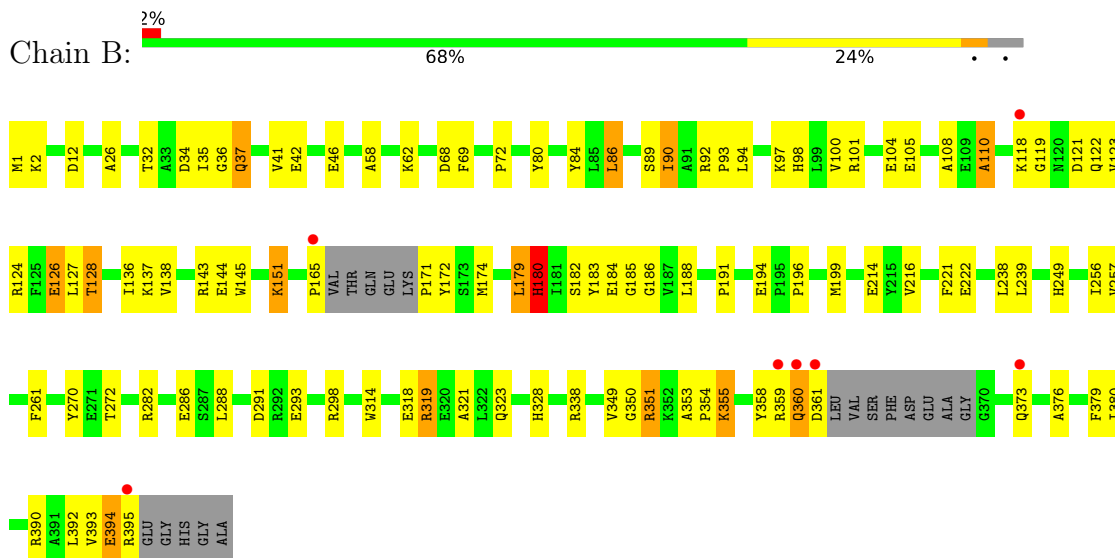
3 Residue-property plots

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

• Molecule 1: Argininosuccinate Synthetase

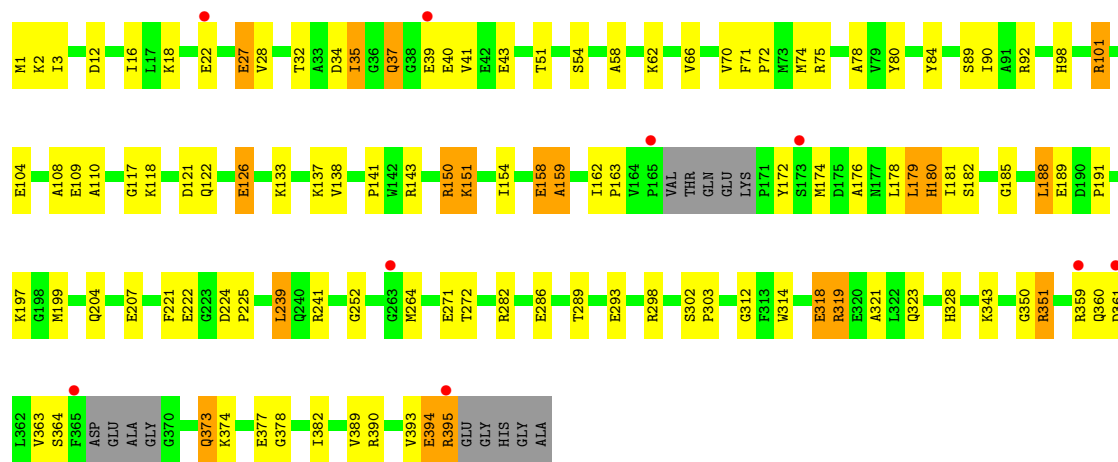


• Molecule 1: Argininosuccinate Synthetase

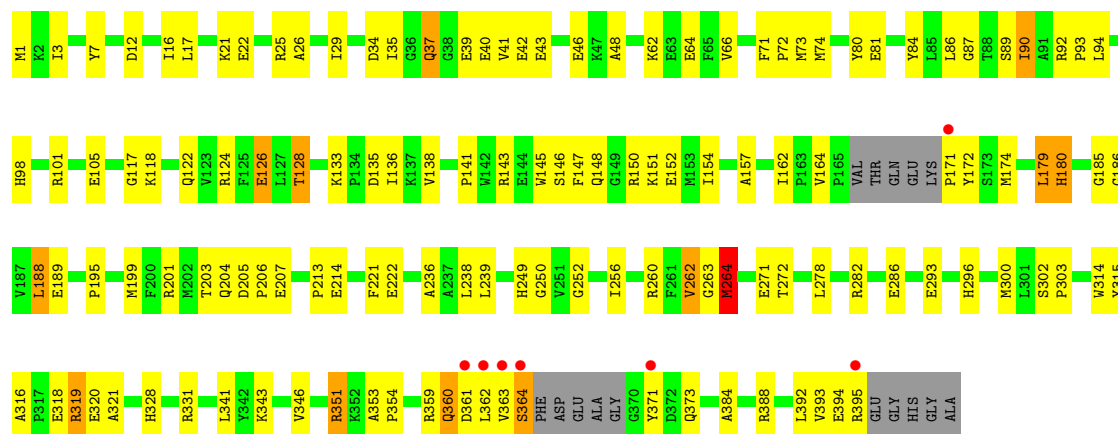


• Molecule 1: Argininosuccinate Synthetase





● 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 10.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.30) 88.0 (10.00-2.30)	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	12404 reflections (9.97%)	wwPDB-VP
Wilson B-factor (Å ²)	32.5	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.45 , 52.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.019 for h,-h-k,-l	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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.46	0/3085	0.98	11/4169 (0.3%)
1	B	0.46	0/3102	0.98	15/4192 (0.4%)
1	C	0.45	0/3135	1.00	13/4237 (0.3%)
1	D	0.45	0/3123	1.01	10/4221 (0.2%)
All	All	0.46	0/12445	0.99	49/16819 (0.3%)

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	221	PHE	CB-CA-C	-16.69	85.42	110.62
1	C	222	GLU	N-CA-C	15.55	132.15	112.24
1	A	222	GLU	N-CA-C	15.16	131.35	112.86
1	D	222	GLU	N-CA-C	14.37	130.76	113.23
1	A	221	PHE	CB-CA-C	-13.91	88.72	110.14
1	C	221	PHE	CB-CA-C	-13.82	87.22	110.16
1	D	221	PHE	CB-CA-C	-13.13	89.13	110.14
1	B	222	GLU	N-CA-C	11.85	127.03	112.58
1	D	145	TRP	N-CA-C	8.31	120.64	110.41
1	A	221	PHE	N-CA-C	-8.23	94.60	108.26
1	C	221	PHE	N-CA-C	-7.89	95.54	108.41
1	D	221	PHE	N-CA-C	-7.73	95.66	108.34
1	A	394	GLU	N-CA-C	-7.69	104.04	113.19
1	C	394	GLU	N-CA-C	-7.53	103.98	113.01
1	D	263	GLY	N-CA-C	7.29	122.32	113.79
1	A	257	VAL	N-CA-C	-6.90	97.15	107.37
1	B	89	SER	N-CA-C	6.74	118.28	111.07
1	C	35	ILE	N-CA-C	-6.55	104.96	111.77
1	A	86	LEU	N-CA-C	6.32	120.27	111.74
1	B	144	GLU	N-CA-C	6.17	120.83	113.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	394	GLU	N-CA-C	-5.95	106.11	113.19
1	D	126	GLU	N-CA-C	5.92	117.53	111.14
1	A	348	VAL	N-CA-C	-5.91	101.92	109.30
1	C	204	GLN	N-CA-C	-5.90	102.75	110.53
1	D	12	ASP	N-CA-C	5.87	117.35	111.07
1	D	256	ILE	N-CA-C	5.82	118.24	108.81
1	C	126	GLU	N-CA-C	5.74	118.48	111.82
1	A	28	VAL	N-CA-C	5.74	116.37	107.99
1	A	89	SER	N-CA-C	5.72	117.19	111.07
1	D	89	SER	N-CA-C	5.71	117.18	111.07
1	C	28	VAL	N-CA-C	5.65	116.23	107.99
1	B	180	HIS	N-CA-C	5.62	115.26	107.73
1	B	126	GLU	N-CA-C	5.57	117.43	111.36
1	C	27	GLU	N-CA-C	-5.52	100.81	109.25
1	B	257	VAL	N-CA-C	-5.46	99.28	107.37
1	D	264	MET	N-CA-C	5.40	118.07	109.81
1	B	12	ASP	N-CA-C	5.38	116.83	111.07
1	C	289	THR	N-CA-C	5.29	119.89	113.38
1	B	86	LEU	N-CA-C	5.28	118.86	111.74
1	B	145	TRP	N-CA-C	5.21	116.65	110.19
1	C	89	SER	N-CA-C	5.17	116.60	110.97
1	B	256	ILE	N-CA-C	5.16	117.47	108.90
1	C	110	ALA	N-CA-C	5.14	116.80	109.14
1	C	78	ALA	N-CA-C	5.11	117.58	109.96
1	B	119	GLY	N-CA-C	5.08	119.53	112.57
1	B	221	PHE	N-CA-C	-5.06	99.01	108.02
1	A	289	THR	N-CA-C	5.03	119.69	113.50
1	B	110	ALA	N-CA-C	5.01	116.69	109.07
1	A	78	ALA	N-CA-C	5.00	117.46	109.96

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	118	0
1	B	3034	0	3018	104	0
1	C	3066	0	3052	110	0
1	D	3055	0	3043	120	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	10	0
3	B	124	0	0	3	0
3	C	105	0	0	6	0
3	D	97	0	0	3	0
All	All	12650	0	12119	407	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 17.

All (407) 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:H	1:B:151:LYS:HD2	1.08	1.08
1:C:37:GLN:H	1:C:37:GLN:NE2	1.53	1.07
1:D:151:LYS:H	1:D:151:LYS:HD2	1.11	1.07
1:C:151:LYS:H	1:C:151:LYS:HD2	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:D:393:VAL:C	1:D:395:ARG:H	1.75	0.92
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:H	1:C:151:LYS:CD	1.83	0.89
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
1:C:151:LYS:HD2	1:C:151:LYS:N	1.91	0.86
1:B:151:LYS:H	1:B:151:LYS:CD	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:HD2	1:D:151:LYS:N	1.95	0.82
1:C:122:GLN:NE2	1:C:143:ARG:HH11	1.78	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:D:393:VAL:C	1:D:395:ARG:N	2.40	0.79
1:A:151:LYS:HE3	1:A:151:LYS:H	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:C:74:MET:HE2	3:C:1315:HOH:O	1.83	0.77
1:A:41:VAL:HG12	1:A:58:ALA:HB1	1.67	0.76
1:A:293:GLU:HB3	1:C:318:GLU:HG2	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:D:262:VAL:HG12	1:D:264:MET:HE3	1.68	0.75
1:C:351:ARG:HD2	1:C:351:ARG:N	2.01	0.74
1:D:74:MET:HE2	3:D:1235:HOH:O	1.86	0.74
1:A:70:VAL:O	1:A:74:MET:HG3	1.86	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: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:HD2	1:B:351:ARG:N	2.02	0.73
1:B:359:ARG:HB3	1:B:361:ASP:OD1	1.88	0.73
1:A:98:HIS:HD2	1:A:101:ARG:NH1	1.85	0.73
1:D:118:LYS:O	1:D:264:MET:HE1	1.90	0.72
1:C:319:ARG:HH21	1:C:323:GLN:NE2	1.87	0.72
1:A:351:ARG:HD2	1:A:351:ARG:N	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:B:293:GLU:HB3	1:D:318:GLU:HG2	1.71	0.70
1:D:126:GLU:HG2	1:D:138:VAL:HG11	1.73	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:18:LYS:CG	1:C:162:ILE:HD11	2.21	0.70
1:C:151:LYS:HE3	3:C:1356:HOH:O	1.90	0.70
1:C:395:ARG:HH11	1:C:395:ARG:CG	2.04	0.70
1:A:123:VAL:HG22	1:B:380:ILE:CD1	2.20	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:179:LEU:O	1:D:272:THR:HG23	1.93	0.69
1:B:393:VAL:C	1:B:395:ARG:H	2.01	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:C:1:MET:HE2	1:C:3:ILE:HD11	1.77	0.66
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:98:HIS:HD2	1:D:101:ARG:HH12	1.44	0.66
1:D:395:ARG:HB3	1:D:395:ARG:CZ	2.26	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:H	1:A:151:LYS:CE	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:H	1:B:37:GLN:HE21	1.44	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:21:LYS:HD3	3:A:1313:HOH:O	1.99	0.63
1:A:151:LYS:H	1:A:151:LYS:CD	2.12	0.63
1:A:393:VAL:C	1:A:395:ARG:H	2.07	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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:VAL:CG1	1:A:58:ALA:HB1	2.30	0.60
1:A:214:GLU:OE1	1:A:249:HIS:HE1	1.84	0.60
1:B:1:MET:HB3	1:B:26:ALA:HB2	1.82	0.60
1:D:90:ILE:HD11	1:D:179:LEU:HD12	1.83	0.60
1:D:360:GLN:HA	1:D:363:VAL:HG22	1.84	0.60
1:C:16:ILE:HG23	1:C:141:PRO:HG2	1.84	0.59
1:B:97:LYS:NZ	3:B:1116:HOH:O	2.35	0.59
1:D:117:GLY:HA2	1:D:122:GLN:NE2	2.17	0.59
1:A:331:ARG:CB	1:A:331:ARG:HH11	2.16	0.59
1:B:359:ARG:O	1:B:361:ASP:N	2.35	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:A:370:GLY:N	3:A:1105:HOH:O	2.35	0.58
1:B:314:TRP:CE3	1:B:319:ARG:HD2	2.38	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:B:393:VAL:C	1:B:395:ARG:N	2.61	0.57
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:D:174:MET:HE1	1:D:201:ARG:HD2	1.85	0.57
1:A:282:ARG:HD2	3:A:1115:HOH:O	2.04	0.57
1:B:35:ILE:N	1:B:37:GLN:HE22	2.00	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: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:D:148:GLN:H	1:D:152:GLU:CD	2.12	0.56
1:B:214:GLU:OE1	1:B:249:HIS:HE1	1.89	0.56
1:C:395:ARG:HG3	1:C:395:ARG:NH1	2.11	0.56
1:C:32:THR:HG22	1:C:41:VAL:HG13	1.87	0.56
1:A:90:ILE:HD11	1:A:179:LEU:HD12	1.86	0.56
1:C:282:ARG:O	1:C:286:GLU:HG3	2.06	0.56
1:B:42:GLU:O	1:B:46:GLU:HG3	2.06	0.56
1:B:282:ARG:O	1:B:286:GLU:HG3	2.05	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:93:PRO:HA	1:B:128:THR:HG21	1.88	0.56
1:B:122:GLN:NE2	1:B:143:ARG:HH11	2.02	0.56
1:D:21:LYS:O	1:D:25:ARG:HA	2.05	0.56
1:C:150:ARG:O	1:C:154:ILE:HG13	2.06	0.55
1:D:316:ALA:HB1	1:D:318:GLU:OE2	2.05	0.55
1:A:302:SER:HB2	1:A:303:PRO:HD3	1.87	0.55
1:B:318:GLU:CG	1:D:293:GLU:HB2	2.37	0.55
1:C:319:ARG:NH2	1:C:323:GLN:HE21	1.96	0.55
1:A:358:TYR:CE2	1:B:191:PRO:HD3	2.40	0.55
1:D:148:GLN:HB2	1:D:152:GLU:OE2	2.06	0.55
1:B:359:ARG:C	1:B:361:ASP:H	2.15	0.55
1:A:293:GLU:HB2	1:C:318:GLU:HG2	1.84	0.55
1:C:122:GLN:HE21	1:C:143:ARG:NH1	2.02	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:66:VAL:HG21	1:A:236:ALA:HA	1.90	0.54
1:A:372:ASP:HB2	1:A:373:GLN:HE21	1.70	0.54
1:B:118:LYS:H	1:B:118:LYS:HD2	1.72	0.54
1:C:343:LYS:HD2	1:D:343:LYS:CD	2.37	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:A:393:VAL:C	1:A:395:ARG:N	2.66	0.53
1:D:174:MET:HE1	1:D:201:ARG:CD	2.37	0.53
1:B:249:HIS:HD2	3:B:1427:HOH:O	1.89	0.53
1:B:379:PHE:HB2	1:C:382:ILE:HG13	1.90	0.53
1:C:70:VAL:O	1:C:74:MET:HG3	2.08	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:O	1:D:262:VAL:CG1	2.55	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: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:D:296:HIS:O	1:D:300:MET:HE3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:180:HIS:ND1	1:C:180:HIS:C	2.67	0.53
1:A:146:SER:O	1:A:147:PHE:C	2.53	0.52
1:C:176:ALA:CB	1:C:181:ILE:HG12	2.39	0.52
1:B:180:HIS:ND1	1:B:180:HIS:C	2.68	0.52
1:A:24:TYR:OH	1:A:144:GLU:OE2	2.28	0.52
1:B:37:GLN:HE21	1:B:37:GLN:N	2.07	0.52
1:B:390:ARG:O	1:B:394:GLU:HG2	2.10	0.52
1:C:373:GLN:H	1:C:373:GLN:NE2	2.07	0.52
1:B:98:HIS:HD2	1:B:101:ARG:NH1	2.07	0.52
1:B:124:ARG:O	1:B:128:THR:HG23	2.10	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:B:80:TYR:HB3	1:B:84:TYR:HB3	1.91	0.52
1:D:150:ARG:HG3	1:D:154:ILE:HD11	1.92	0.52
1:D:87:GLY:HA2	1:D:179:LEU:HD13	1.91	0.51
1:C:37:GLN:HB2	1:C:39:GLU:HG2	1.92	0.51
1:C:394:GLU:O	1:C:395:ARG:HB2	2.10	0.51
1:B:86:LEU:O	1:B:90:ILE:HG13	2.11	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:C:109:GLU:O	1:C:137:LYS:HE2	2.11	0.51
1:D:393:VAL:O	1:D:395:ARG:N	2.44	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:B:359:ARG:C	1:B:361:ASP:N	2.68	0.51
1:C:92:ARG:HE	1:C:121:ASP:CG	2.19	0.51
1:D:93:PRO:HA	1:D:128:THR:HG21	1.93	0.51
1:D:146:SER:O	1:D:147:PHE:C	2.53	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:B:179:LEU:HD22	1:B:272:THR:HG23	1.93	0.50
1:C:373:GLN:HG2	1:C:374:LYS:N	2.27	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: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:D:360:GLN:O	1:D:364:SER:HB2	2.11	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:373:GLN:H	1:A:373:GLN:CD	2.16	0.49
1:D:351:ARG:HD2	1:D:351:ARG:H	1.74	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:C:390:ARG:O	1:C:394:GLU:HG2	2.13	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
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: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:A:84:TYR:CE2	1:A:86:LEU:HA	2.49	0.48
1:A:339:LEU:HD12	1:A:339:LEU:N	2.29	0.48
1:B:36:GLY:N	1:B:37:GLN:HE21	2.12	0.48
1:C:178:LEU:HB3	1:C:239:LEU:HD13	1.96	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:73:MET:HE3	3:A:1221:HOH:O	2.13	0.47
1:A:350:GLY:C	1:A:351:ARG:HD2	2.38	0.47
1:C:174:MET:HA	1:C:182:SER:O	2.14	0.47
1:D:262:VAL:CG1	1:D:264:MET:HE3	2.42	0.47
1:A:179:LEU:HD22	1:A:272:THR:HG23	1.97	0.47
1:D:314:TRP:CE3	1:D:319:ARG:HD2	2.49	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: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:C:395:ARG:CG	1:C:395:ARG:NH1	2.68	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:B:185:GLY:O	1:B:188:LEU:HB2	2.15	0.47
1:D:80:TYR:CE2	1:D:81:GLU:HG3	2.49	0.47
1:A:3:ILE:HD13	1:A:139:ILE:HD12	1.97	0.47
1:C:1:MET:HE2	1:C:3:ILE:CD1	2.45	0.47
1:D:22:GLU:HA	1:D:25:ARG:NH1	2.30	0.47
1:C:75:ARG:NH2	3:C:1156:HOH:O	2.48	0.46

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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	0.46
1:C:389:VAL:O	1:C:393:VAL:HG23	2.15	0.46
1:A:154:ILE:HG22	1:A:155:ALA:N	2.31	0.46
1:B:238:LEU:HD23	1:B:238:LEU:C	2.40	0.46
1:C:101:ARG:NH2	3:C:1140:HOH:O	2.34	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:328:HIS:CD2	1:D:321:ALA:HA	2.50	0.46
1:C:314:TRP:O	1:C:319:ARG:HD3	2.15	0.46
1:D:84:TYR:CZ	1:D:86:LEU:HA	2.50	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:B:298:ARG:HD2	1:B:298:ARG:C	2.41	0.46
1:D:203:THR:HB	1:D:250:GLY:HA2	1.98	0.46
1:B:350:GLY:C	1:B:351:ARG:HD2	2.41	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:D:238:LEU:C	1:D:238:LEU:HD23	2.41	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:HD12	1:D:29:ILE:N	2.31	0.45
1:A:298:ARG:C	1:A:298:ARG:HD2	2.41	0.45
1:C:374:LYS:HE2	1:C:377:GLU:OE2	2.17	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:80:TYR:HB3	1:A:84:TYR:HB3	1.98	0.44
1:A:151:LYS:HE3	1:A:151:LYS:N	2.25	0.44
1:A:331:ARG:CB	1:A:331:ARG:NH1	2.79	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:B:69:PHE:C	1:B:72:PRO:HD2	2.42	0.44
1:B:174:MET:HG2	1:B:183:TYR:CD2	2.53	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:A:100:VAL:O	1:A:104:GLU:HG3	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:PRO:HD3	1:B:358:TYR:CE2	2.52	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: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:B:126:GLU:CG	1:B:138:VAL:HG11	2.42	0.44
1:A:257:VAL:HG11	1:B:288:LEU:HD23	2.00	0.44
1:B:321:ALA:HA	1:D:328:HIS:CD2	2.53	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: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:H	1:A:197:LYS:HG2	1.57	0.43
1:C:117:GLY:HA3	1:D:373:GLN:OE1	2.18	0.43
1:C:312:GLY:HA2	3:C:1083:HOH:O	2.19	0.43
1:D:7:TYR:CD2	1:D:48:ALA:HB2	2.53	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:316:ALA:O	1:D:320:GLU:HG3	2.19	0.43
1:D:341:LEU:HD22	1:D:346:VAL:HG22	2.00	0.43
1:A:17:LEU:HD23	1:A:17:LEU:C	2.44	0.43
1:B:122:GLN:NE2	1:B:143:ARG:HD3	2.33	0.43
1:A:122:GLN:OE1	1:A:143:ARG:HD3	2.19	0.43
1:C:109:GLU:HG2	1:C:137:LYS:CE	2.49	0.43
1:C:37:GLN:NE2	1:C:37:GLN:N	2.33	0.43
1:A:185:GLY:HA2	1:A:189:GLU:HG3	2.01	0.43
1:D:73:MET:CE	1:D:90:ILE:HG12	2.48	0.43
1:A:4:VAL:O	1:A:111:ILE:HA	2.19	0.43
1:A:24:TYR:O	1:A:25:ARG:C	2.61	0.43
1:A:179:LEU:O	1:A:272:THR:HG23	2.18	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: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:D:92:ARG:HB2	1:D:93:PRO:HD3	2.01	0.42
1:A:174:MET:HA	1:A:182:SER:O	2.19	0.42
1:A:321:ALA:HA	1:C:328:HIS:CD2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:158:GLU:O	1:C:159:ALA:C	2.62	0.42
1:C:172:TYR:CZ	1:C:199:MET:HG3	2.54	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:D:180:HIS:ND1	1:D:180:HIS:C	2.77	0.42
1:B:238:LEU:HD23	1:B:238:LEU:O	2.19	0.42
1:A:7:TYR:CE2	1:A:9:GLY:HA2	2.54	0.42
1:A:15:ILE:HG23	1:A:156:TYR:CD2	2.55	0.42
1:A:85:LEU:HD23	1:A:85:LEU:HA	1.83	0.42
1:D:205:ASP:OD2	1:D:207:GLU:HB2	2.20	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
1:B:136:ILE:HG13	1:B:137:LYS:N	2.34	0.41
1:D:17:LEU:HD23	1:D:17:LEU:C	2.44	0.41
1:D:34:ASP:HB2	1:D:41:VAL:HG11	2.02	0.41
1:D:171:PRO:HG2	1:D:172:TYR:H	1.84	0.41
1:A:181:ILE:HD12	1:A:202:MET:HB2	2.02	0.41
1:B:351:ARG:N	1:B:351:ARG:CD	2.80	0.41
1:C:80:TYR:HB3	1:C:84:TYR:HB3	2.02	0.41
1:C:162:ILE:O	1:C:163:PRO:C	2.62	0.41
1:C:343:LYS:HD2	1:D:343:LYS:HD2	2.02	0.41
1:D:359:ARG:NH2	1:D:361:ASP:OD1	2.52	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:A:345:ASN:HB3	1:A:347:TYR:CE1	2.56	0.41
1:C:40:GLU:O	1:C:43:GLU:N	2.50	0.41
1:D:360:GLN:O	1:D:364:SER:N	2.53	0.41
1:A:282:ARG:O	1:A:286:GLU:HG3	2.21	0.41
1:B:68:ASP:O	1:B:72:PRO:HG2	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: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:80:TYR:CD2	1:A:81:GLU:HG3	2.56	0.41
1:D:1:MET:HE2	1:D:3:ILE:CG1	2.51	0.41
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:180:HIS:ND1	1:A:180:HIS:C	2.78	0.40
1:A:185:GLY:O	1:A:188:LEU:HB2	2.21	0.40
1:C:314:TRP:O	1:C:319:ARG:CD	2.69	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:359:ARG:NE	1:C:361:ASP:OD1	2.54	0.40
1:C:71:PHE:HB2	1:C:72:PRO:HD3	2.03	0.40
1:C:207:GLU:OE1	1:D:213:PRO:HG2	2.21	0.40
1:C:298:ARG:C	1:C:298:ARG:HD2	2.46	0.40
1:C:393:VAL:C	1:C:395:ARG:H	2.28	0.40
1:A:118:LYS:H	1:A:118:LYS:HD2	1.87	0.40
1:A:180:HIS:HB2	1:A:271:GLU:O	2.22	0.40
1:B:179:LEU:HD22	1:B:272:THR:CG2	2.50	0.40
1:D:74:MET:HE1	1:D:278:LEU:HD12	2.01	0.40
1:D:195:PRO:HB3	1:D:199:MET:CE	2.51	0.40
1:D:205:ASP:O	1:D:206:PRO:C	2.64	0.40
1:D:359:ARG:NH2	3:D:1096:HOH:O	2.53	0.40
1:A:75:ARG:NH2	3:A:1346:HOH:O	2.54	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:C:151:LYS:CD	1:C:151:LYS:N	2.58	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%)	36	46
1	B	376/400 (94%)	362 (96%)	13 (4%)	1 (0%)	36	46
1	C	380/400 (95%)	365 (96%)	14 (4%)	1 (0%)	36	46
1	D	379/400 (95%)	365 (96%)	14 (4%)	0	100	100
All	All	1509/1600 (94%)	1452 (96%)	54 (4%)	3 (0%)	43	55

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 ⓘ

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%)	290 (95%)	14 (5%)	24	36
1	B	306/319 (96%)	289 (94%)	17 (6%)	19	28
1	C	310/319 (97%)	292 (94%)	18 (6%)	18	26
1	D	309/319 (97%)	291 (94%)	18 (6%)	18	26
All	All	1229/1276 (96%)	1162 (94%)	67 (6%)	19	28

All (67) 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	151	LYS
1	A	154	ILE
1	A	179	LEU
1	A	180	HIS
1	A	239	LEU
1	A	318	GLU
1	A	319	ARG
1	A	349	VAL
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	121	ASP

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Mol	Chain	Res	Type
1	B	128	THR
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	291	ASP
1	B	319	ARG
1	B	351	ARG
1	B	355	LYS
1	C	22	GLU
1	C	37	GLN
1	C	62	LYS
1	C	90	ILE
1	C	101	ARG
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	318	GLU
1	C	319	ARG
1	C	351	ARG
1	C	373	GLN
1	C	395	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	179	LEU
1	D	180	HIS
1	D	188	LEU
1	D	239	LEU
1	D	262	VAL
1	D	264	MET

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Mol	Chain	Res	Type
1	D	319	ARG
1	D	351	ARG
1	D	360	GLN
1	D	364	SER
1	D	371	TYR
1	D	394	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) 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	180	HIS
1	C	323	GLN
1	C	328	HIS
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	204	GLN
1	D	249	HIS
1	D	323	GLN

5.3.3 RNA ⓘ

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](#)

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	C	2510	-	4,4,4	0.47	0	6,6,6	0.07	0
2	SO4	A	510	-	4,4,4	0.35	0	6,6,6	0.18	0
2	SO4	B	1510	-	4,4,4	0.36	0	6,6,6	0.21	0
2	SO4	D	3510	-	4,4,4	0.43	0	6,6,6	0.23	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 ⓘ

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%)	-0.49	5 (1%) 75 76	20, 31, 54, 70	0
1	B	382/400 (95%)	-0.56	7 (1%) 67 69	22, 31, 48, 83	0
1	C	386/400 (96%)	-0.34	9 (2%) 61 63	23, 34, 60, 82	0
1	D	385/400 (96%)	-0.35	7 (1%) 67 69	20, 34, 59, 82	0
All	All	1533/1600 (95%)	-0.43	28 (1%) 67 69	20, 32, 56, 83	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	364	SER	4.8
1	A	395	ARG	4.1
1	C	365	PHE	3.9
1	C	165	PRO	3.8
1	D	371	TYR	3.5
1	D	395	ARG	3.4
1	C	395	ARG	3.3
1	D	171	PRO	3.2
1	C	39	GLU	3.0
1	D	363	VAL	2.9
1	C	361	ASP	2.8
1	B	360	GLN	2.8
1	B	373	GLN	2.7
1	B	361	ASP	2.7
1	A	171	PRO	2.7
1	D	361	ASP	2.7
1	B	395	ARG	2.7
1	A	359	ARG	2.4
1	D	362	LEU	2.4
1	C	22	GLU	2.4
1	B	165	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	197	LYS	2.3
1	B	359	ARG	2.3
1	C	263	GLY	2.2
1	C	359	ARG	2.1
1	B	118	LYS	2.0
1	A	149	GLY	2.0
1	C	173	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	D	3510	5/5	0.96	0.08	45,49,51,54	0
2	SO4	C	2510	5/5	0.97	0.08	50,51,53,55	0
2	SO4	A	510	5/5	0.97	0.08	43,44,46,46	0
2	SO4	B	1510	5/5	0.99	0.04	41,44,47,49	0

6.5 Other polymers [i](#)

There are no such residues in this entry.