



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:37 AM UTC

PDB ID : 2IP4 / pdb_00002ip4
Title : Crystal Structure of Glycinamide Ribonucleotide Synthetase from *Thermus thermophilus* HB8
Authors : Sampei, G.; Baba, S.; Kanagawa, M.; Yanai, H.; Ishii, T.; Kawai, H.; Fukai, Y.; Ebihara, A.; Nakagawa, N.; Kawai, G.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2006-10-11
Resolution : 2.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 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

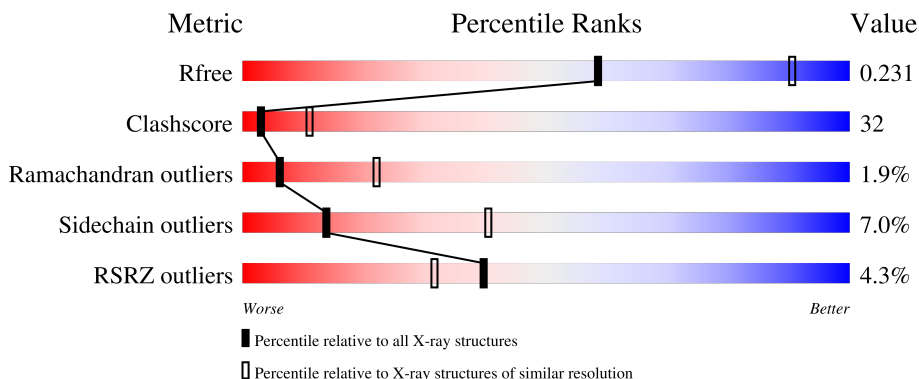
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	417	 4% 51% 42% 6%
1	B	417	 4% 53% 40% 6%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6268 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 Phosphoribosylamine--glycine ligase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	414	Total 3112	C 1979	N 554	O 570	S 1	Se 8	0	0	0
1	B	414	Total 3112	C 1979	N 554	O 570	S 1	Se 8	0	0	0

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

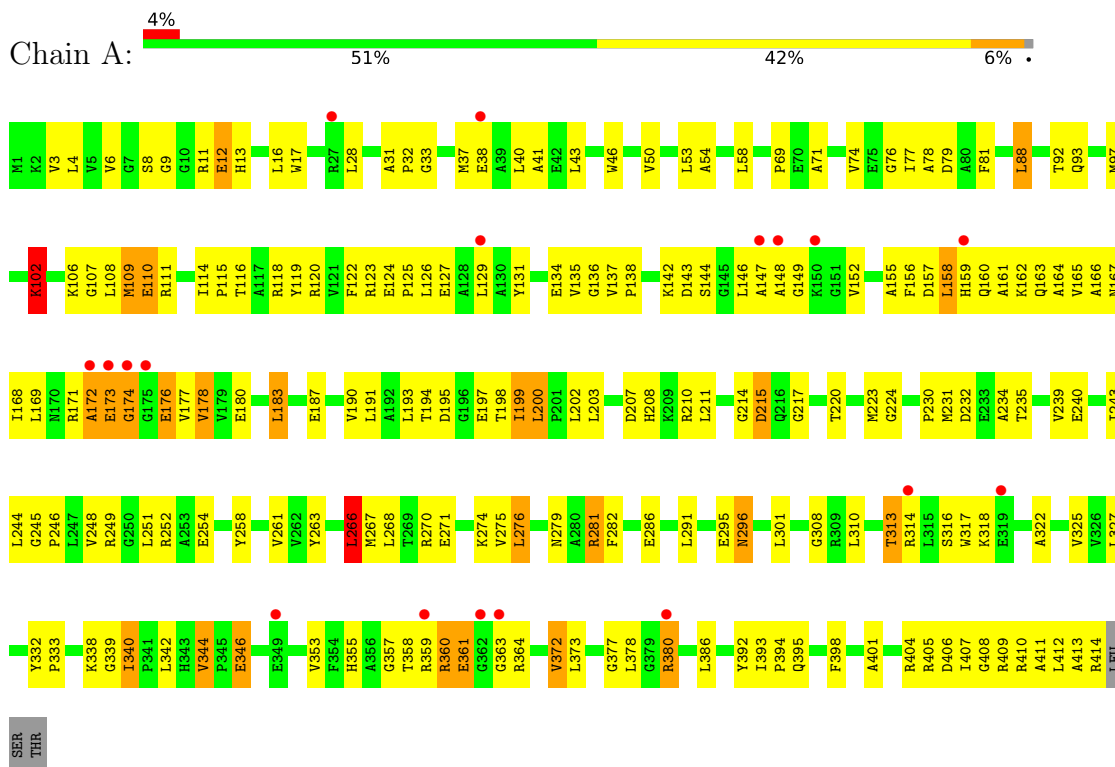
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total 11	O 11	0	0
3	B	13	Total 13	O 13	0	0

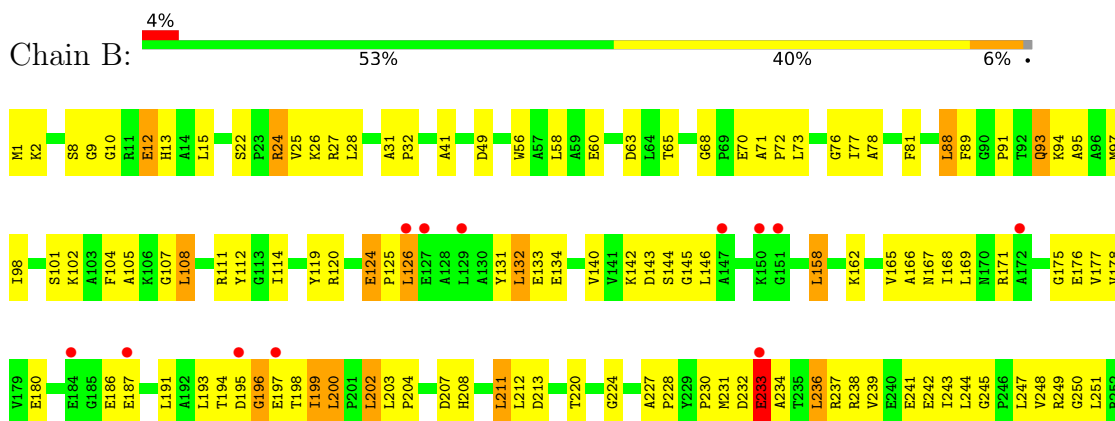
3 Residue-property plots [i](#)

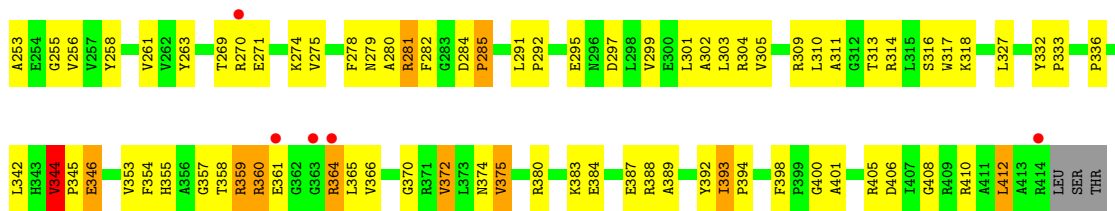
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Phosphoribosylamine--glycine ligase



- Molecule 1: Phosphoribosylamine--glycine ligase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.31Å 98.07Å 121.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.00 – 2.80 60.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (60.00-2.80) 99.3 (60.00-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.77Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.218 , 0.238 0.228 , 0.231	Depositor DCC
R_{free} test set	1117 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	36.4	Xtrriage
Anisotropy	0.312	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6268	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 41.76 % 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 2.2514e-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, 0.333 respectively for untwinned datasets, and 0.375, 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: 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	0.44	0/3165	1.00	13/4285 (0.3%)
1	B	0.45	0/3165	1.01	10/4285 (0.2%)
All	All	0.44	0/6330	1.01	23/8570 (0.3%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	158	LEU	N-CA-C	-12.88	96.92	112.89
1	B	196	GLY	N-CA-C	-9.83	101.99	114.37
1	A	291	LEU	CA-C-N	-8.06	111.50	119.64
1	A	291	LEU	C-N-CA	-8.06	111.50	119.64
1	A	176	GLU	N-CA-C	7.80	120.65	110.43
1	A	380	ARG	N-CA-C	-7.61	104.05	112.72
1	B	158	LEU	N-CA-C	-7.48	103.15	111.82
1	A	372	VAL	N-CA-C	6.75	117.36	111.56
1	B	124	GLU	CA-C-N	6.63	126.88	119.32
1	B	124	GLU	C-N-CA	6.63	126.88	119.32
1	B	393	ILE	CB-CA-C	-6.62	107.41	113.70
1	B	372	VAL	N-CA-C	6.12	117.48	111.67
1	A	124	GLU	CA-C-N	6.01	127.35	119.84
1	A	124	GLU	C-N-CA	6.01	127.35	119.84
1	A	102	LYS	N-CA-C	-5.98	104.70	112.23
1	B	375	VAL	N-CA-C	-5.97	99.81	108.17
1	A	215	ASP	N-CA-C	-5.76	103.44	111.39
1	B	212	LEU	N-CA-C	5.68	118.83	110.59
1	B	344	VAL	CA-C-N	5.31	126.13	120.13
1	B	344	VAL	C-N-CA	5.31	126.13	120.13
1	A	276	LEU	N-CA-C	-5.26	106.92	113.55
1	A	266	LEU	N-CA-C	5.14	116.65	108.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	156	PHE	N-CA-C	-5.08	106.33	112.88

There are no chirality outliers.

There are no planarity outliers.

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	3112	0	3174	205	0
1	B	3112	0	3174	195	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
3	A	11	0	0	6	0
3	B	13	0	0	3	0
All	All	6268	0	6348	400	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 32.

All (400) 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:360:ARG:HH11	1:A:360:ARG:HB2	1.23	1.04
1:A:193:LEU:HD13	1:A:202:LEU:HD21	1.36	1.03
1:A:16:LEU:HD22	1:A:37:MSE:HE2	1.41	1.02
1:A:267:MSE:HG2	1:A:276:LEU:HD11	1.38	1.02
1:B:108:LEU:HB3	1:B:114:ILE:HD12	1.43	1.00
1:B:126:LEU:H	1:B:126:LEU:HD12	1.31	0.94
1:A:157:ASP:HB3	1:A:160:GLN:H	1.31	0.93
1:A:263:TYR:HB3	1:A:279:ASN:HB2	1.55	0.89
1:B:94:LYS:HG2	3:B:429:HOH:O	1.75	0.87
1:B:231:MSE:HE3	1:B:236:LEU:HB2	1.58	0.85
1:A:108:LEU:HB3	1:A:114:ILE:HG13	1.58	0.84
1:A:279:ASN:ND2	1:A:281:ARG:HH21	1.75	0.84
1:A:360:ARG:HB2	1:A:360:ARG:NH1	1.93	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:ILE:HG12	1:A:248:VAL:HG21	1.62	0.81
1:A:144:SER:HB3	1:A:176:GLU:HB3	1.61	0.81
1:A:296:ASN:HD22	1:A:296:ASN:H	1.25	0.81
1:A:279:ASN:HD22	1:A:281:ARG:HH21	1.28	0.81
1:A:207:ASP:HB2	1:A:286:GLU:OE1	1.82	0.80
1:B:15:LEU:HD23	1:B:291:LEU:HD11	1.64	0.79
1:A:243:ILE:HD12	1:A:275:VAL:HG23	1.64	0.78
1:B:13:HIS:HD2	1:B:355:HIS:H	1.31	0.78
1:B:104:PHE:O	1:B:108:LEU:HD22	1.84	0.77
1:B:65:THR:HB	1:B:88:LEU:HB2	1.66	0.77
1:B:1:MSE:HA	1:B:26:LYS:HE2	1.65	0.77
1:A:183:LEU:H	1:A:183:LEU:HD12	1.49	0.76
1:B:200:LEU:N	1:B:200:LEU:HD12	2.01	0.75
1:A:115:PRO:HG2	1:A:243:ILE:HD11	1.67	0.75
1:B:112:TYR:HE1	1:B:249:ARG:HD2	1.53	0.74
1:A:183:LEU:HD12	1:A:183:LEU:N	2.02	0.73
1:A:9:GLY:O	1:A:12:GLU:HG2	1.88	0.73
1:A:395:GLN:HE21	1:A:395:GLN:HA	1.53	0.73
1:A:210:ARG:NH1	1:A:217:GLY:O	2.21	0.73
1:A:224:GLY:HA3	1:A:407:ILE:HD13	1.71	0.71
1:B:126:LEU:H	1:B:126:LEU:CD1	2.04	0.71
1:B:9:GLY:H	1:B:12:GLU:CG	2.04	0.71
1:B:358:THR:HB	1:B:365:LEU:HD11	1.71	0.71
1:B:263:TYR:HB3	1:B:279:ASN:HB2	1.73	0.70
1:B:22:SER:HB2	1:B:297:ASP:OD1	1.91	0.70
1:B:120:ARG:HD3	1:B:131:TYR:CZ	2.27	0.69
1:B:186:GLU:OE2	1:B:410:ARG:HG3	1.92	0.69
1:B:208:HIS:HD2	1:B:224:GLY:O	1.75	0.69
1:B:28:LEU:O	1:B:41:ALA:HB1	1.93	0.69
1:B:13:HIS:CD2	1:B:355:HIS:H	2.09	0.69
1:A:135:VAL:HG23	1:A:136:GLY:N	2.08	0.69
1:B:193:LEU:CD1	1:B:202:LEU:HD21	2.24	0.68
1:A:168:ILE:HD12	1:A:177:VAL:HG22	1.74	0.68
1:B:94:LYS:HD2	1:B:256:VAL:HG12	1.73	0.68
1:B:88:LEU:C	1:B:88:LEU:HD12	2.19	0.67
1:A:6:VAL:O	1:A:69:PRO:HD2	1.95	0.67
1:A:93:GLN:HG2	1:A:97:MSE:HE2	1.77	0.67
1:A:296:ASN:HD21	1:A:316:SER:H	1.39	0.67
1:A:183:LEU:HD23	1:A:274:LYS:HD2	1.76	0.67
1:A:120:ARG:HB2	1:A:131:TYR:CE1	2.30	0.67
1:B:360:ARG:HH11	1:B:360:ARG:HB3	1.60	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:PRO:HD3	1:B:169:LEU:HD23	1.76	0.67
1:A:210:ARG:O	1:A:404:ARG:HD2	1.95	0.66
1:B:193:LEU:HD23	1:B:305:VAL:HG21	1.78	0.66
1:B:295:GLU:OE1	1:B:318:LYS:HE3	1.94	0.66
1:B:108:LEU:CB	1:B:114:ILE:HD12	2.21	0.66
1:A:13:HIS:HD2	1:A:355:HIS:H	1.42	0.65
1:A:187:GLU:HB2	1:A:267:MSE:HE2	1.77	0.65
1:B:108:LEU:HB3	1:B:114:ILE:CD1	2.22	0.65
1:A:106:LYS:NZ	1:A:110:GLU:HG3	2.11	0.65
1:B:107:GLY:O	1:B:111:ARG:HB2	1.97	0.65
1:B:375:VAL:HG21	1:B:393:ILE:HD13	1.77	0.65
1:A:119:TYR:O	1:A:120:ARG:HG3	1.97	0.65
1:B:56:TRP:CZ2	1:B:60:GLU:HG3	2.31	0.65
1:B:112:TYR:CE1	1:B:249:ARG:HD2	2.32	0.64
1:A:190:VAL:HG12	1:A:190:VAL:O	1.97	0.64
1:B:220:THR:HB	1:B:332:TYR:OH	1.98	0.64
1:A:50:VAL:CG1	1:A:77:ILE:HA	2.27	0.64
1:A:239:VAL:O	1:A:243:ILE:HB	1.99	0.63
1:A:144:SER:C	1:A:146:LEU:H	2.07	0.63
1:A:106:LYS:HZ1	1:A:110:GLU:HG3	1.64	0.63
1:B:104:PHE:CD2	1:B:108:LEU:HD21	2.32	0.63
1:A:358:THR:O	1:A:359:ARG:HD2	1.98	0.62
1:B:104:PHE:CE2	1:B:108:LEU:HD21	2.34	0.62
1:A:74:VAL:O	1:A:93:GLN:HG3	1.97	0.62
1:B:357:GLY:O	1:B:372:VAL:HG13	1.99	0.62
1:B:231:MSE:HE3	1:B:236:LEU:CB	2.29	0.62
1:B:364:ARG:HH11	1:B:364:ARG:HG2	1.65	0.62
1:A:231:MSE:HE2	1:A:268:LEU:HD11	1.82	0.62
1:A:122:PHE:CD1	1:A:127:GLU:HB3	2.35	0.61
1:A:165:VAL:HG13	1:A:166:ALA:H	1.65	0.61
1:A:142:LYS:HB3	1:A:178:VAL:HG13	1.81	0.61
1:A:393:ILE:N	1:A:394:PRO:HD2	2.16	0.61
1:A:249:ARG:NH1	1:A:249:ARG:HB3	2.14	0.61
1:A:332:TYR:CD1	1:A:333:PRO:HA	2.35	0.61
1:A:168:ILE:HA	1:A:171:ARG:HH12	1.66	0.60
1:A:296:ASN:HD22	1:A:296:ASN:N	1.98	0.60
1:A:143:ASP:HA	1:A:168:ILE:HD13	1.83	0.60
1:A:327:LEU:O	1:A:398:PHE:HE1	1.85	0.60
1:B:102:LYS:O	1:B:105:ALA:HB3	2.02	0.60
1:A:114:ILE:HD11	1:A:246:PRO:HB2	1.84	0.60
1:B:126:LEU:HD12	1:B:126:LEU:N	2.11	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:142:LYS:O	1:B:178:VAL:HG12	2.02	0.59
1:A:198:THR:HG22	1:A:200:LEU:HD13	1.84	0.59
1:A:372:VAL:HG12	1:A:373:LEU:HG	1.82	0.59
1:B:9:GLY:H	1:B:12:GLU:HG3	1.66	0.59
1:B:398:PHE:CG	1:B:401:ALA:HB2	2.38	0.59
1:B:213:ASP:HA	1:B:406:ASP:OD1	2.03	0.59
1:A:223:MSE:O	1:A:407:ILE:HD11	2.02	0.59
1:B:389:ALA:O	1:B:393:ILE:HG12	2.02	0.59
1:A:266:LEU:HA	1:A:276:LEU:HG	1.85	0.58
1:B:28:LEU:C	1:B:41:ALA:HB1	2.28	0.58
1:B:71:ALA:HB3	1:B:72:PRO:HD3	1.84	0.58
1:A:392:TYR:C	1:A:394:PRO:HD2	2.27	0.58
1:B:203:LEU:HG	1:B:317:TRP:CD2	2.37	0.58
1:B:94:LYS:O	1:B:98:ILE:HG12	2.03	0.58
1:B:295:GLU:HB2	1:B:316:SER:HB2	1.84	0.58
1:A:261:VAL:HB	1:A:281:ARG:HA	1.85	0.58
1:A:342:LEU:HD12	1:A:373:LEU:HD11	1.86	0.58
1:A:135:VAL:HG23	1:A:136:GLY:H	1.67	0.58
1:A:157:ASP:CB	1:A:160:GLN:H	2.13	0.57
1:B:142:LYS:HB3	1:B:178:VAL:CG1	2.34	0.57
1:A:171:ARG:HB2	1:A:171:ARG:NH1	2.20	0.57
1:B:193:LEU:HD23	1:B:305:VAL:CG2	2.34	0.57
1:A:50:VAL:HG13	1:A:77:ILE:HA	1.87	0.57
1:B:13:HIS:HD2	1:B:355:HIS:N	2.02	0.57
1:B:166:ALA:C	1:B:168:ILE:H	2.13	0.57
1:A:195:ASP:HB2	1:A:308:GLY:HA2	1.87	0.57
1:A:325:VAL:HG13	1:A:407:ILE:HD12	1.87	0.56
1:B:393:ILE:N	1:B:394:PRO:HD2	2.21	0.56
1:A:157:ASP:HB3	1:A:160:GLN:N	2.13	0.56
1:A:109:MSE:HE3	1:A:116:THR:OG1	2.05	0.56
1:A:313:THR:HG22	1:A:314:ARG:H	1.70	0.56
1:B:380:ARG:H	1:B:384:GLU:CD	2.14	0.56
1:B:342:LEU:HD11	1:B:372:VAL:HG11	1.88	0.56
1:A:137:VAL:HA	1:A:138:PRO:C	2.30	0.56
1:B:193:LEU:HD12	1:B:202:LEU:HD21	1.88	0.56
1:B:342:LEU:HD11	1:B:372:VAL:CG1	2.36	0.55
1:B:270:ARG:HH11	1:B:270:ARG:HG2	1.72	0.55
1:A:53:LEU:HD22	1:A:81:PHE:HZ	1.71	0.55
1:B:204:PRO:HG3	1:B:231:MSE:HE2	1.88	0.55
1:A:168:ILE:HA	1:A:171:ARG:NH1	2.22	0.55
1:A:407:ILE:HG13	3:A:420:HOH:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:ARG:HH12	1:A:281:ARG:HH11	1.54	0.55
1:A:38:GLU:HG2	1:A:43:LEU:HD22	1.88	0.55
1:A:171:ARG:HB2	1:A:171:ARG:HH11	1.72	0.55
1:B:332:TYR:CD1	1:B:333:PRO:HA	2.42	0.55
1:A:171:ARG:O	1:A:172:ALA:HB2	2.05	0.55
1:A:281:ARG:C	1:A:281:ARG:HD2	2.31	0.55
1:A:398:PHE:CG	1:A:401:ALA:HB2	2.41	0.55
1:A:296:ASN:ND2	1:A:316:SER:H	2.05	0.55
1:A:161:ALA:O	1:A:164:ALA:HB3	2.07	0.55
1:A:263:TYR:HE2	1:A:286:GLU:HG3	1.73	0.55
1:B:12:GLU:OE2	1:B:68:GLY:HA3	2.07	0.54
1:A:9:GLY:H	1:A:12:GLU:CG	2.19	0.54
1:B:250:GLY:O	1:B:253:ALA:HB3	2.08	0.54
1:B:304:ARG:HD2	1:B:309:ARG:O	2.08	0.54
1:B:119:TYR:C	1:B:120:ARG:HG3	2.33	0.54
1:B:345:PRO:O	1:B:346:GLU:C	2.51	0.54
1:B:200:LEU:HD13	1:B:310:LEU:HD21	1.90	0.54
1:A:410:ARG:HG2	1:A:410:ARG:HH11	1.73	0.54
1:B:2:LYS:HG2	1:B:27:ARG:HB2	1.89	0.54
1:B:200:LEU:N	1:B:200:LEU:CD1	2.69	0.53
1:A:295:GLU:HB2	1:A:316:SER:HB2	1.91	0.53
1:B:344:VAL:HG21	1:B:353:VAL:HG11	1.90	0.53
1:A:165:VAL:HG13	1:A:166:ALA:N	2.23	0.53
1:B:124:GLU:OE2	1:B:125:PRO:HD2	2.07	0.53
1:A:243:ILE:HD12	1:A:275:VAL:CG2	2.37	0.53
1:A:313:THR:C	1:A:314:ARG:HD2	2.34	0.53
1:A:168:ILE:CD1	1:A:177:VAL:HG22	2.36	0.53
1:B:194:THR:HA	1:B:198:THR:O	2.09	0.53
1:A:210:ARG:HB2	1:A:215:ASP:HA	1.89	0.53
1:B:24:ARG:HH11	1:B:24:ARG:CB	2.22	0.53
1:A:267:MSE:CG	1:A:276:LEU:HD11	2.26	0.53
1:B:309:ARG:HB2	1:B:309:ARG:NH1	2.24	0.53
1:B:9:GLY:O	1:B:12:GLU:HG2	2.09	0.52
1:B:15:LEU:CD2	1:B:291:LEU:HD11	2.36	0.52
1:A:249:ARG:HH11	1:A:249:ARG:CB	2.22	0.52
1:B:360:ARG:HB3	1:B:360:ARG:NH1	2.25	0.52
1:B:197:GLU:HG3	1:B:198:THR:N	2.24	0.52
1:B:233:GLU:HA	1:B:236:LEU:HB3	1.92	0.52
1:A:414:ARG:HA	1:A:414:ARG:NE	2.24	0.52
1:A:224:GLY:HA3	1:A:407:ILE:CD1	2.39	0.52
1:B:94:LYS:HG3	1:B:95:ALA:N	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:TYR:O	1:B:120:ARG:HG3	2.09	0.52
1:A:4:LEU:HD21	1:A:53:LEU:HD21	1.92	0.52
1:B:195:ASP:HB2	1:B:310:LEU:HB3	1.90	0.52
1:A:232:ASP:OD2	1:A:234:ALA:HB3	2.11	0.51
1:B:359:ARG:HB3	1:B:366:VAL:HG22	1.91	0.51
1:B:1:MSE:HB3	1:B:63:ASP:OD2	2.09	0.51
1:A:118:ARG:NH2	1:A:135:VAL:HA	2.26	0.51
1:B:144:SER:C	1:B:146:LEU:H	2.18	0.51
1:B:93:GLN:O	1:B:97:MSE:HG3	2.11	0.51
1:B:125:PRO:CD	1:B:169:LEU:HD23	2.40	0.51
1:B:392:TYR:C	1:B:394:PRO:HD2	2.36	0.51
1:A:173:GLU:O	1:A:174:GLY:O	2.28	0.51
1:B:1:MSE:HE1	1:B:299:VAL:HG11	1.91	0.51
1:B:199:ILE:C	1:B:200:LEU:HD12	2.36	0.51
1:A:169:LEU:HD12	1:A:169:LEU:N	2.26	0.51
1:A:270:ARG:HB2	1:A:271:GLU:OE2	2.11	0.51
1:A:361:GLU:OE2	1:A:364:ARG:HB2	2.11	0.51
1:A:407:ILE:N	3:A:420:HOH:O	2.31	0.51
1:B:198:THR:HG22	1:B:199:ILE:N	2.25	0.51
1:B:70:GLU:O	1:B:73:LEU:HB2	2.11	0.51
1:B:243:ILE:HD13	1:B:275:VAL:HG23	1.93	0.50
1:A:202:LEU:HB3	1:A:317:TRP:CH2	2.46	0.50
1:A:37:MSE:SE	3:A:430:HOH:O	2.79	0.50
1:A:77:ILE:HG23	1:A:78:ALA:N	2.25	0.50
1:B:383:LYS:O	1:B:387:GLU:HG3	2.12	0.50
1:A:406:ASP:HB2	1:A:409:ARG:HG3	1.94	0.50
1:A:11:ARG:NH2	1:A:69:PRO:HA	2.27	0.50
1:B:13:HIS:CD2	1:B:354:PHE:HA	2.47	0.50
1:B:227:ALA:HA	1:B:228:PRO:C	2.36	0.50
1:A:77:ILE:CG2	1:A:78:ALA:N	2.75	0.50
1:B:89:PHE:CE2	1:B:261:VAL:HG23	2.47	0.50
1:B:195:ASP:HB3	1:B:198:THR:H	1.76	0.50
1:A:169:LEU:N	1:A:169:LEU:CD1	2.74	0.50
1:A:17:TRP:HA	1:A:40:LEU:HD13	1.94	0.49
1:A:249:ARG:HB3	1:A:249:ARG:HH11	1.75	0.49
1:B:202:LEU:HB3	1:B:317:TRP:CH2	2.47	0.49
1:A:135:VAL:CG2	1:A:136:GLY:N	2.74	0.49
1:B:8:SER:N	1:B:12:GLU:HG3	2.27	0.49
1:A:16:LEU:CD2	1:A:37:MSE:HE2	2.29	0.49
1:A:395:GLN:HA	1:A:395:GLN:NE2	2.25	0.49
1:B:1:MSE:HE1	1:B:299:VAL:CG1	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:GLY:N	1:B:12:GLU:HG2	2.28	0.49
1:B:236:LEU:O	1:B:239:VAL:HG22	2.13	0.49
1:B:88:LEU:HD12	1:B:89:PHE:N	2.27	0.49
1:B:360:ARG:HH11	1:B:360:ARG:CB	2.26	0.49
1:B:233:GLU:CD	1:B:233:GLU:H	2.21	0.49
1:A:157:ASP:HB2	1:A:160:GLN:HB3	1.94	0.49
1:B:89:PHE:CZ	1:B:261:VAL:HG23	2.47	0.49
1:B:251:LEU:HD13	1:B:258:TYR:HB3	1.95	0.49
1:B:125:PRO:HG2	1:B:126:LEU:HD12	1.94	0.48
1:A:146:LEU:O	1:A:148:ALA:N	2.46	0.48
1:A:410:ARG:HG2	1:A:410:ARG:NH1	2.28	0.48
1:B:9:GLY:N	1:B:12:GLU:CG	2.75	0.48
1:A:158:LEU:HD23	1:A:158:LEU:O	2.13	0.48
1:B:93:GLN:O	1:B:94:LYS:C	2.57	0.48
1:B:28:LEU:HB2	1:B:41:ALA:HB2	1.95	0.48
1:B:162:LYS:O	1:B:165:VAL:HG12	2.13	0.48
1:B:220:THR:HB	1:B:332:TYR:CZ	2.49	0.48
1:B:9:GLY:H	1:B:12:GLU:HG2	1.77	0.48
1:B:237:ARG:HG3	1:B:237:ARG:HH11	1.79	0.47
1:A:171:ARG:O	1:A:172:ALA:CB	2.62	0.47
1:A:183:LEU:H	1:A:183:LEU:CD1	2.24	0.47
1:B:1:MSE:CA	1:B:26:LYS:HE2	2.40	0.47
1:A:12:GLU:CD	1:A:12:GLU:H	2.22	0.47
1:A:120:ARG:HD3	1:A:131:TYR:CE2	2.49	0.47
1:B:76:GLY:HA2	3:B:428:HOH:O	2.13	0.47
1:B:199:ILE:HD11	1:B:245:GLY:HA2	1.96	0.47
1:A:194:THR:HB	1:A:199:ILE:HG23	1.96	0.47
1:A:413:ALA:O	1:A:414:ARG:HB2	2.14	0.47
1:B:25:VAL:HG11	1:B:28:LEU:HD21	1.97	0.47
1:B:208:HIS:CE1	1:B:408:GLY:H	2.33	0.47
1:A:50:VAL:HG11	1:A:77:ILE:HA	1.96	0.47
1:A:214:GLY:O	1:A:215:ASP:C	2.58	0.47
1:A:266:LEU:HB3	1:A:275:VAL:HA	1.97	0.47
1:A:314:ARG:HD2	1:A:314:ARG:N	2.30	0.47
1:A:412:LEU:C	1:A:414:ARG:H	2.23	0.47
1:B:98:ILE:O	1:B:101:SER:HB3	2.15	0.47
1:A:8:SER:N	1:A:12:GLU:HG3	2.30	0.47
1:B:91:PRO:HG3	1:B:280:ALA:HB1	1.97	0.47
1:A:144:SER:C	1:A:146:LEU:N	2.72	0.46
1:A:249:ARG:O	1:A:252:ARG:HG2	2.15	0.46
1:B:24:ARG:HB2	1:B:24:ARG:NH1	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:258:TYR:CD1	1:B:258:TYR:C	2.94	0.46
1:B:364:ARG:HH11	1:B:364:ARG:CG	2.28	0.46
1:A:135:VAL:CG2	1:A:136:GLY:H	2.28	0.46
1:B:251:LEU:HD13	1:B:258:TYR:CB	2.45	0.46
1:B:269:THR:HG21	1:B:274:LYS:HD2	1.97	0.46
1:A:360:ARG:HD2	1:A:363:GLY:C	2.41	0.46
1:A:211:LEU:HD13	1:A:211:LEU:O	2.16	0.46
1:A:106:LYS:HD3	1:A:107:GLY:N	2.31	0.46
1:B:253:ALA:C	1:B:255:GLY:H	2.24	0.46
1:B:256:VAL:HG23	1:B:256:VAL:O	2.15	0.46
1:B:327:LEU:HD21	1:B:393:ILE:HD12	1.96	0.46
1:A:119:TYR:HB2	1:A:180:GLU:HB3	1.97	0.46
1:A:296:ASN:H	1:A:296:ASN:ND2	2.02	0.46
1:A:235:THR:O	1:A:239:VAL:HG23	2.16	0.46
1:A:76:GLY:HA2	1:A:93:GLN:HB2	1.96	0.45
1:A:144:SER:CB	1:A:176:GLU:HB3	2.41	0.45
1:A:211:LEU:HD23	1:A:220:THR:HG21	1.98	0.45
1:A:142:LYS:HG3	1:A:152:VAL:HG22	1.97	0.45
1:B:238:ARG:O	1:B:242:GLU:HB2	2.16	0.45
1:A:338:LYS:HA	3:A:426:HOH:O	2.16	0.45
1:A:386:LEU:HD13	1:A:412:LEU:HD11	1.99	0.45
1:A:414:ARG:HA	1:A:414:ARG:CZ	2.45	0.45
1:B:88:LEU:HD12	1:B:89:PHE:C	2.42	0.45
1:A:251:LEU:HD21	1:A:258:TYR:CD2	2.51	0.45
1:B:15:LEU:HD23	1:B:291:LEU:CD1	2.42	0.45
1:B:364:ARG:N	1:B:364:ARG:HD2	2.31	0.45
1:B:346:GLU:CD	1:B:346:GLU:H	2.24	0.45
1:B:144:SER:O	1:B:146:LEU:N	2.50	0.45
1:B:244:LEU:O	1:B:247:LEU:HB3	2.17	0.45
1:A:122:PHE:HD1	1:A:127:GLU:HB3	1.82	0.45
1:A:200:LEU:HD22	1:A:310:LEU:HG	2.00	0.45
1:A:240:GLU:HA	1:A:244:LEU:HB2	1.98	0.45
1:B:133:GLU:HG2	1:B:158:LEU:HD21	1.98	0.45
1:B:302:ALA:O	1:B:303:LEU:C	2.60	0.45
1:A:79:ASP:OD1	1:A:92:THR:HB	2.17	0.44
1:A:108:LEU:HB3	1:A:114:ILE:CG1	2.39	0.44
1:A:149:GLY:O	1:A:152:VAL:HG23	2.17	0.44
1:A:342:LEU:CD1	1:A:373:LEU:HD11	2.47	0.44
1:B:336:PRO:HB3	1:B:370:GLY:HA2	1.99	0.44
1:A:33:GLY:HA2	1:A:37:MSE:SE	2.67	0.44
1:A:203:LEU:HG	1:A:317:TRP:CD2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:THR:HG22	1:A:314:ARG:N	2.33	0.44
1:B:88:LEU:C	1:B:88:LEU:CD1	2.89	0.44
1:B:193:LEU:O	1:B:199:ILE:HA	2.17	0.44
1:A:11:ARG:NH1	1:A:281:ARG:HH11	2.16	0.44
1:B:354:PHE:HB2	1:B:374:ASN:HB2	1.99	0.44
1:A:102:LYS:HD2	1:A:119:TYR:CZ	2.53	0.44
1:A:197:GLU:CD	1:A:252:ARG:HH22	2.26	0.44
1:A:208:HIS:CE1	1:A:408:GLY:H	2.35	0.44
1:B:31:ALA:HA	1:B:32:PRO:HA	1.89	0.44
1:B:232:ASP:O	1:B:234:ALA:N	2.51	0.44
1:A:123:ARG:HG2	1:A:176:GLU:HG3	1.99	0.44
1:A:16:LEU:HD21	1:A:41:ALA:HB3	1.99	0.44
1:B:94:LYS:HG3	1:B:95:ALA:H	1.81	0.44
1:B:143:ASP:HB2	1:B:168:ILE:HD13	2.00	0.44
1:A:411:ALA:O	1:A:412:LEU:C	2.61	0.44
1:B:384:GLU:OE2	1:B:388:ARG:NH2	2.47	0.44
1:A:71:ALA:HA	1:A:74:VAL:HG22	2.00	0.44
1:A:187:GLU:HB2	1:A:267:MSE:CE	2.48	0.44
1:B:207:ASP:OD1	1:B:208:HIS:N	2.49	0.43
1:B:358:THR:HB	1:B:365:LEU:CD1	2.45	0.43
1:A:46:TRP:CD2	1:A:53:LEU:HD12	2.53	0.43
1:A:16:LEU:HB2	1:A:37:MSE:HE3	1.99	0.43
1:B:243:ILE:HD13	1:B:275:VAL:CG2	2.47	0.43
1:A:109:MSE:O	1:A:109:MSE:SE	2.86	0.43
1:B:120:ARG:HD3	1:B:131:TYR:CE1	2.53	0.43
1:B:275:VAL:HG11	1:B:278:PHE:CE2	2.54	0.43
1:A:281:ARG:HD2	1:A:281:ARG:O	2.18	0.43
1:B:22:SER:OG	1:B:24:ARG:HG2	2.19	0.43
1:B:211:LEU:C	1:B:211:LEU:CD2	2.91	0.43
1:B:234:ALA:O	1:B:237:ARG:HB3	2.18	0.43
1:A:157:ASP:CG	1:A:159:HIS:HB3	2.43	0.43
1:B:165:VAL:HG13	1:B:166:ALA:N	2.34	0.43
1:B:291:LEU:N	1:B:292:PRO:HD2	2.33	0.43
1:B:77:ILE:HG23	1:B:78:ALA:N	2.34	0.42
1:B:105:ALA:HA	1:B:108:LEU:HD23	2.00	0.42
1:B:166:ALA:C	1:B:168:ILE:N	2.77	0.42
1:A:183:LEU:N	1:A:183:LEU:CD1	2.74	0.42
1:B:199:ILE:HD13	1:B:248:VAL:CG2	2.50	0.42
1:A:12:GLU:CD	1:A:12:GLU:N	2.77	0.42
1:A:78:ALA:HA	1:A:88:LEU:HG	2.01	0.42
1:A:393:ILE:N	1:A:394:PRO:CD	2.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MSE:CE	1:B:299:VAL:HG11	2.50	0.42
1:A:211:LEU:HA	1:A:223:MSE:HG3	2.01	0.42
1:B:196:GLY:HA2	1:B:258:TYR:O	2.20	0.42
1:A:339:GLY:O	1:A:340:ILE:C	2.61	0.42
1:A:138:PRO:HA	1:A:155:ALA:O	2.19	0.42
1:A:346:GLU:H	1:A:346:GLU:CD	2.28	0.42
1:B:10:GLY:HA3	1:B:374:ASN:OD1	2.20	0.42
1:A:191:LEU:HD23	1:A:263:TYR:HA	2.00	0.42
1:B:191:LEU:HD12	3:B:421:HOH:O	2.19	0.42
1:A:211:LEU:HD13	1:A:211:LEU:C	2.44	0.41
1:A:404:ARG:C	3:A:420:HOH:O	2.63	0.41
1:B:400:GLY:O	1:B:401:ALA:C	2.63	0.41
1:B:131:TYR:O	1:B:134:GLU:HB3	2.20	0.41
1:B:313:THR:HG22	1:B:314:ARG:N	2.35	0.41
1:A:296:ASN:N	1:A:296:ASN:ND2	2.65	0.41
1:B:49:ASP:OD1	1:B:49:ASP:C	2.62	0.41
1:B:237:ARG:O	1:B:241:GLU:HG2	2.20	0.41
1:B:412:LEU:HD12	1:B:412:LEU:HA	1.87	0.41
1:A:377:GLY:C	1:A:378:LEU:HD12	2.45	0.41
1:A:202:LEU:HB3	1:A:317:TRP:HH2	1.85	0.41
1:B:208:HIS:CD2	1:B:224:GLY:O	2.64	0.41
1:B:271:GLU:HB2	1:B:274:LYS:NZ	2.36	0.41
1:B:281:ARG:HD3	1:B:282:PHE:O	2.20	0.41
1:A:107:GLY:HA3	1:A:111:ARG:HH21	1.85	0.41
1:A:125:PRO:O	1:A:126:LEU:C	2.64	0.41
1:A:163:GLN:O	1:A:167:ASN:HB2	2.21	0.41
1:A:261:VAL:HG11	1:A:282:PHE:N	2.35	0.41
1:A:322:ALA:HA	1:A:377:GLY:O	2.21	0.41
1:B:187:GLU:O	1:B:187:GLU:HG3	2.20	0.41
1:A:318:LYS:HB3	1:A:380:ARG:HH22	1.85	0.41
1:B:304:ARG:HA	1:B:304:ARG:HD3	1.94	0.41
1:A:3:VAL:HB	1:A:28:LEU:HD23	2.03	0.41
1:A:160:GLN:O	1:A:160:GLN:HG2	2.21	0.41
1:A:168:ILE:CA	1:A:171:ARG:HH12	2.32	0.41
1:A:344:VAL:HG21	1:A:353:VAL:HG11	2.02	0.41
1:A:386:LEU:CD1	1:A:412:LEU:HD11	2.51	0.41
1:B:140:VAL:HB	1:B:180:GLU:HG3	2.03	0.41
1:B:301:LEU:O	1:B:304:ARG:HB2	2.21	0.41
1:B:358:THR:CB	1:B:365:LEU:HD11	2.46	0.41
1:A:31:ALA:HA	1:A:32:PRO:HA	1.74	0.41
1:A:129:LEU:HD22	1:A:162:LYS:HG2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:SER:HA	3:A:430:HOH:O	2.21	0.40
1:B:81:PHE:CG	1:B:88:LEU:HD23	2.56	0.40
1:B:132:LEU:HD12	1:B:132:LEU:HA	1.77	0.40
1:B:165:VAL:HG23	1:B:177:VAL:HG11	2.03	0.40
1:B:197:GLU:HG3	1:B:198:THR:OG1	2.21	0.40
1:A:162:LYS:C	1:A:164:ALA:N	2.78	0.40
1:B:1:MSE:H2	1:B:26:LYS:HG2	1.85	0.40
1:B:232:ASP:C	1:B:234:ALA:N	2.78	0.40
1:B:281:ARG:C	1:B:281:ARG:CD	2.95	0.40
1:A:53:LEU:O	1:A:54:ALA:C	2.64	0.40
1:A:245:GLY:O	1:A:249:ARG:HG3	2.21	0.40
1:B:144:SER:HB3	1:B:176:GLU:HB3	2.04	0.40
1:A:131:TYR:O	1:A:134:GLU:HB3	2.22	0.40
1:B:197:GLU:CG	1:B:198:THR:N	2.83	0.40
1:A:78:ALA:HB3	1:A:92:THR:HA	2.03	0.40
1:A:243:ILE:CD1	1:A:275:VAL:HG23	2.45	0.40
1:B:284:ASP:HA	1:B:285:PRO:HA	1.69	0.40
1:B:336:PRO:CB	1:B:370:GLY:HA2	2.52	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	412/417 (99%)	365 (89%)	40 (10%)	7 (2%)	7	25
1	B	412/417 (99%)	375 (91%)	28 (7%)	9 (2%)	5	19
All	All	824/834 (99%)	740 (90%)	68 (8%)	16 (2%)	6	22

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	147	ALA
1	A	174	GLY
1	B	171	ARG
1	A	172	ALA
1	A	173	GLU
1	B	167	ASN
1	B	233	GLU
1	B	311	ALA
1	A	230	PRO
1	B	93	GLN
1	B	230	PRO
1	B	361	GLU
1	A	357	GLY
1	B	145	GLY
1	A	340	ILE
1	B	175	GLY

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	308/303 (102%)	287 (93%)	21 (7%)	14	42
1	B	308/303 (102%)	286 (93%)	22 (7%)	13	39
All	All	616/606 (102%)	573 (93%)	43 (7%)	14	40

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	GLU
1	A	58	LEU
1	A	88	LEU
1	A	102	LYS
1	A	109	MSE
1	A	110	GLU
1	A	178	VAL
1	A	183	LEU

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Mol	Chain	Res	Type
1	A	199	ILE
1	A	200	LEU
1	A	254	GLU
1	A	266	LEU
1	A	281	ARG
1	A	296	ASN
1	A	301	LEU
1	A	313	THR
1	A	344	VAL
1	A	346	GLU
1	A	360	ARG
1	A	361	GLU
1	A	405	ARG
1	B	12	GLU
1	B	24	ARG
1	B	58	LEU
1	B	88	LEU
1	B	108	LEU
1	B	126	LEU
1	B	132	LEU
1	B	199	ILE
1	B	200	LEU
1	B	202	LEU
1	B	211	LEU
1	B	233	GLU
1	B	236	LEU
1	B	281	ARG
1	B	285	PRO
1	B	344	VAL
1	B	346	GLU
1	B	359	ARG
1	B	360	ARG
1	B	364	ARG
1	B	405	ARG
1	B	412	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	HIS
1	A	21	GLN
1	A	82	GLN

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Mol	Chain	Res	Type
1	A	208	HIS
1	A	216	GLN
1	A	279	ASN
1	A	296	ASN
1	A	343	HIS
1	A	395	GLN
1	B	13	HIS
1	B	21	GLN
1	B	93	GLN
1	B	159	HIS
1	B	167	ASN
1	B	208	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	B	419	-	4,4,4	0.42	0	6,6,6	0.13	0
2	SO4	B	418	-	4,4,4	0.40	0	6,6,6	0.18	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	418	-	4,4,4	0.41	0	6,6,6	0.16	0
2	SO4	A	419	-	4,4,4	0.35	0	6,6,6	0.08	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	406/417 (97%)	0.12	18 (4%) 39 31	3, 28, 76, 147	0
1	B	406/417 (97%)	-0.00	17 (4%) 40 32	3, 24, 80, 136	0
All	All	812/834 (97%)	0.06	35 (4%) 40 31	3, 26, 80, 147	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	126	LEU	5.1
1	A	150	LYS	4.5
1	A	148	ALA	4.2
1	B	129	LEU	3.9
1	B	127	GLU	3.7
1	A	380	ARG	3.6
1	A	363	GLY	3.4
1	A	159	HIS	3.1
1	B	147	ALA	2.9
1	A	362	GLY	2.8
1	B	233	GLU	2.8
1	A	173	GLU	2.8
1	A	349	GLU	2.8
1	B	172	ALA	2.8
1	B	361	GLU	2.8
1	A	172	ALA	2.8
1	B	414	ARG	2.7
1	A	27	ARG	2.7
1	A	38	GLU	2.6
1	B	184	GLU	2.6
1	A	314	ARG	2.6
1	A	147	ALA	2.5
1	B	364	ARG	2.5
1	A	175	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	363	GLY	2.4
1	B	270	ARG	2.4
1	B	150	LYS	2.4
1	B	197	GLU	2.3
1	A	319	GLU	2.3
1	A	129	LEU	2.3
1	B	151	GLY	2.2
1	A	174	GLY	2.2
1	B	187	GLU	2.2
1	A	359	ARG	2.1
1	B	195	ASP	2.1

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(Å ²)	Q<0.9
2	SO4	A	419	5/5	0.92	0.11	49,49,51,52	0
2	SO4	B	418	5/5	0.93	0.11	48,49,50,50	0
2	SO4	B	419	5/5	0.94	0.12	29,30,32,33	0
2	SO4	A	418	5/5	0.96	0.06	33,33,33,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.