Full wwPDB X-ray Structure Validation Report (i)

Jun 13, 2024 – 12:51 AM EDT

PDB ID	:	3KD9
Title	:	Crystal structure of pyridine nucleotide disulfide oxidoreductase from Pyrococcus horikoshii
Authors	:	Agarwal, R.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on	:	2009-10-22
Resolution	:	2.75 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 (i) 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 (i)) were used in the production of this report:

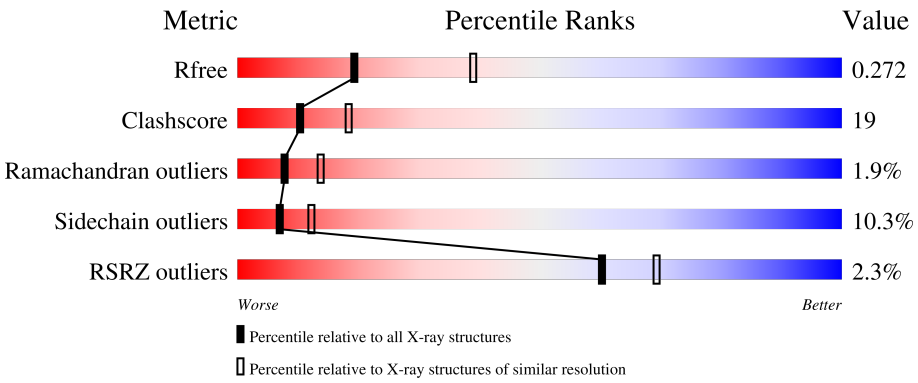
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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}	130704	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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 <=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	449	2% 58% 32% • 7%
1	B	449	2% 61% 26% 6% 7%
1	C	449	2% 62% 27% • 7%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 9568 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 Coenzyme A disulfide reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	418	3144	2004	543	587	10	0	0	0
1	B	418	3151	2010	543	588	10	0	0	0
1	C	418	3148	2009	543	586	10	0	0	0

There are 33 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP O58308
A	2	SER	-	expression tag	UNP O58308
A	3	LEU	-	expression tag	UNP O58308
A	442	GLU	-	expression tag	UNP O58308
A	443	GLY	-	expression tag	UNP O58308
A	444	HIS	-	expression tag	UNP O58308
A	445	HIS	-	expression tag	UNP O58308
A	446	HIS	-	expression tag	UNP O58308
A	447	HIS	-	expression tag	UNP O58308
A	448	HIS	-	expression tag	UNP O58308
A	449	HIS	-	expression tag	UNP O58308
B	1	MET	-	expression tag	UNP O58308
B	2	SER	-	expression tag	UNP O58308
B	3	LEU	-	expression tag	UNP O58308
B	442	GLU	-	expression tag	UNP O58308
B	443	GLY	-	expression tag	UNP O58308
B	444	HIS	-	expression tag	UNP O58308
B	445	HIS	-	expression tag	UNP O58308
B	446	HIS	-	expression tag	UNP O58308
B	447	HIS	-	expression tag	UNP O58308
B	448	HIS	-	expression tag	UNP O58308
B	449	HIS	-	expression tag	UNP O58308
C	1	MET	-	expression tag	UNP O58308

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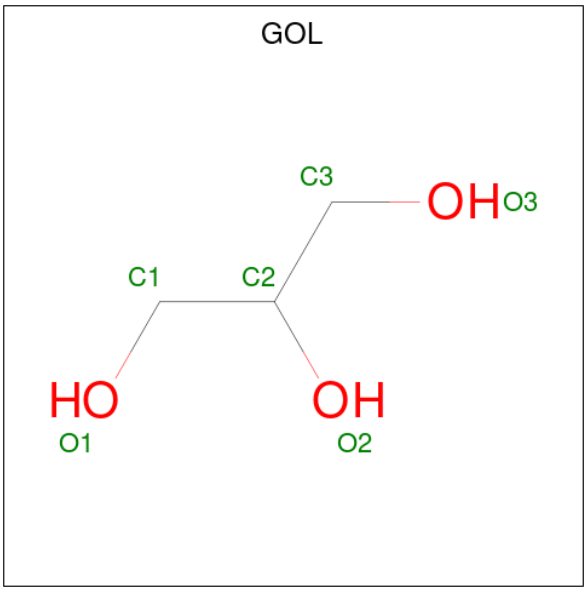
Chain	Residue	Modelled	Actual	Comment	Reference
C	2	SER	-	expression tag	UNP O58308
C	3	LEU	-	expression tag	UNP O58308
C	442	GLU	-	expression tag	UNP O58308
C	443	GLY	-	expression tag	UNP O58308
C	444	HIS	-	expression tag	UNP O58308
C	445	HIS	-	expression tag	UNP O58308
C	446	HIS	-	expression tag	UNP O58308
C	447	HIS	-	expression tag	UNP O58308
C	448	HIS	-	expression tag	UNP O58308
C	449	HIS	-	expression tag	UNP O58308

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	6	3	3	0	0

- Molecule 3 is water.

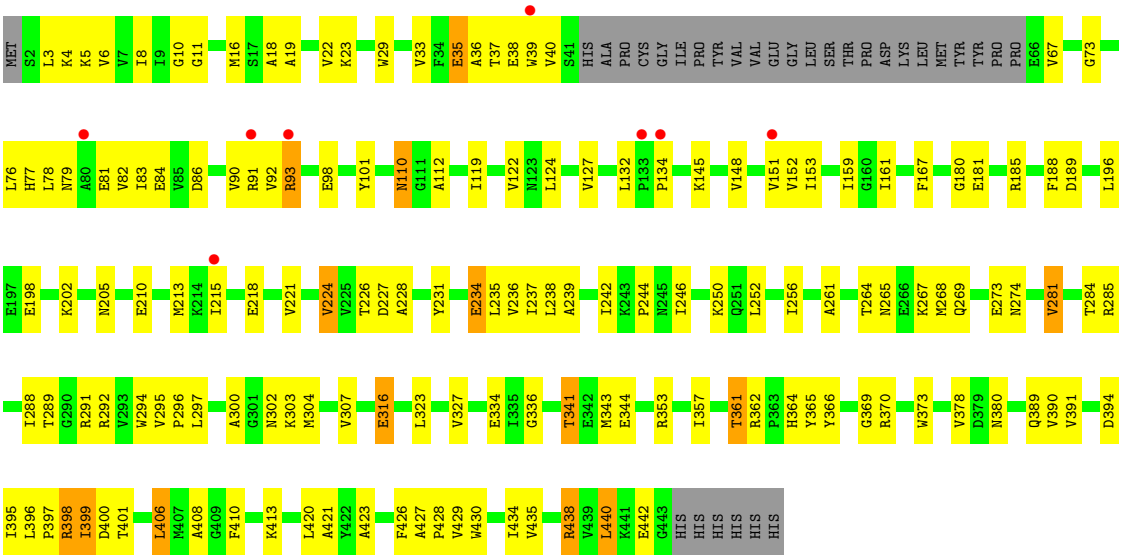
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	42	42	42	0	0
			Total	O		
3	B	32	32	32	0	0
			Total	O		
3	C	45	45	45	0	0

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.

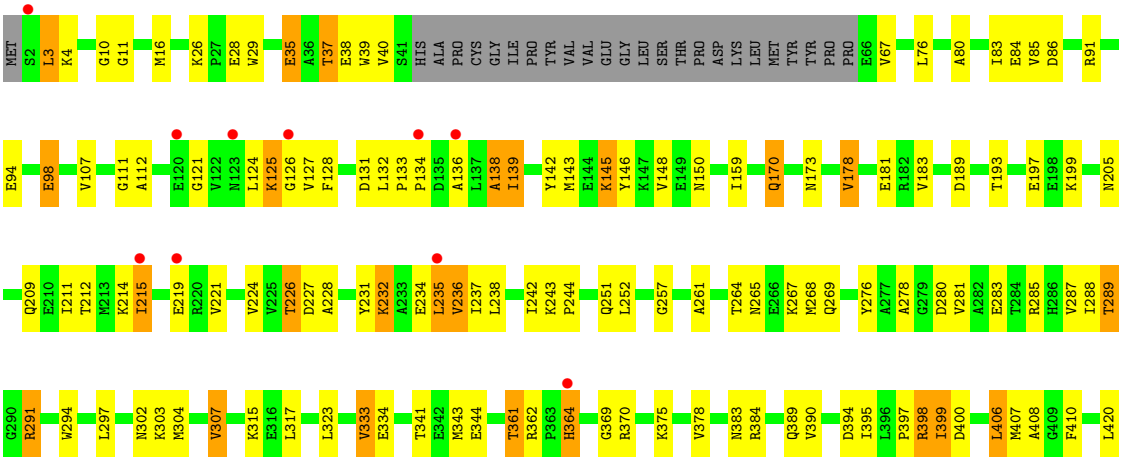
• Molecule 1: Coenzyme A disulfide reductase

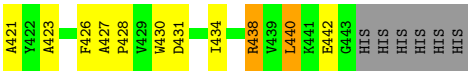
Chain A:



• Molecule 1: Coenzyme A disulfide reductase

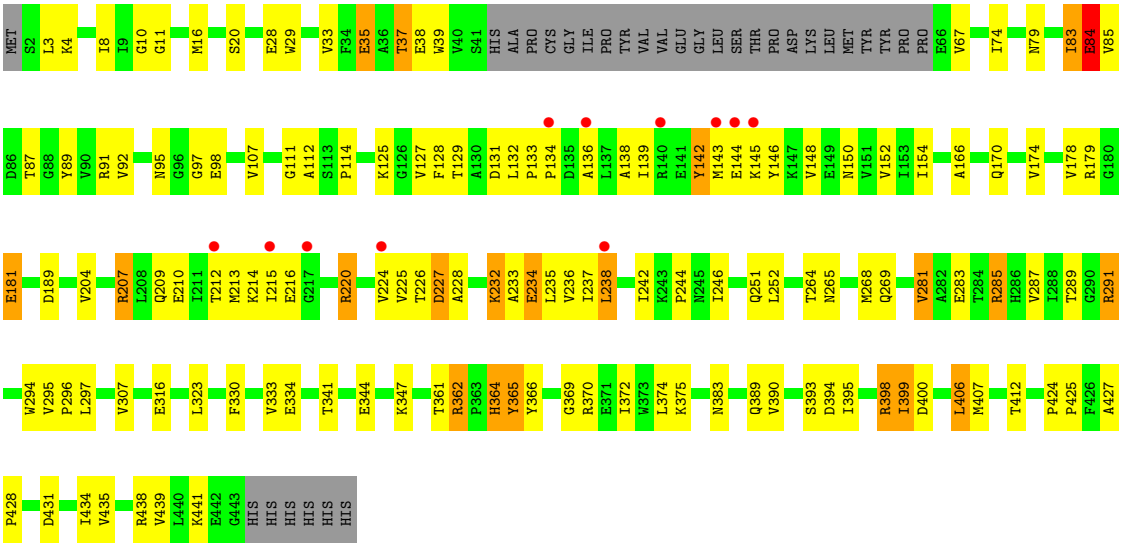
Chain B:





• Molecule 1: Coenzyme A disulfide reductase

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	114.81Å 114.81Å 353.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.31 – 2.75 49.31 – 2.75	Depositor EDS
% Data completeness (in resolution range)	96.3 (49.31-2.75) 96.3 (49.31-2.75)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.59 (at 2.77Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.244 , 0.284 0.236 , 0.272	Depositor DCC
R_{free} test set	1827 reflections (2.94%)	wwPDB-VP
Wilson B-factor (Å^2)	57.0	Xtriage
Anisotropy	0.060	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 33.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	9568	wwPDB-VP
Average B, all atoms (Å^2)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.56% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL

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	0/3201	0.72	1/4342 (0.0%)
1	B	0.48	0/3209	0.71	0/4353
1	C	0.48	0/3206	0.72	0/4349
All	All	0.47	0/9616	0.72	1/13044 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	235	LEU	CA-CB-CG	5.09	127.00	115.30

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	3144	0	3123	131	0
1	B	3151	0	3130	123	0
1	C	3148	0	3128	113	0
2	A	6	0	8	0	0
3	A	42	0	0	2	0
3	B	32	0	0	1	0
3	C	45	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	9568	0	9389	352	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 19.

All (352) 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:37:THR:HG23	1:B:39:TRP:H	1.26	1.00
1:A:93:ARG:HH11	1:A:93:ARG:HB3	1.29	0.98
1:C:37:THR:HG23	1:C:39:TRP:H	1.29	0.96
1:A:399:ILE:HD11	1:B:421:ALA:HB2	1.49	0.92
1:A:11:GLY:HA3	1:A:35:GLU:HG2	1.49	0.91
1:B:189:ASP:H	1:B:389:GLN:HE22	1.17	0.90
1:C:189:ASP:H	1:C:389:GLN:HE22	1.10	0.90
1:A:189:ASP:H	1:A:389:GLN:HE22	0.95	0.90
1:A:110:ASN:H	1:A:110:ASN:HD22	1.23	0.87
1:C:220:ARG:HH11	1:C:220:ARG:HB3	1.38	0.87
1:B:84:GLU:HA	1:B:251:GLN:HE22	1.40	0.86
1:A:265:ASN:HD21	1:A:269:GLN:HE21	1.22	0.85
1:B:361:THR:HG23	1:B:362:ARG:HD3	1.59	0.85
1:C:129:THR:HG22	1:C:238:LEU:HB2	1.59	0.84
1:B:215:ILE:HG22	1:B:224:VAL:HA	1.60	0.84
1:B:127:VAL:HA	1:B:236:VAL:HG23	1.60	0.83
1:A:110:ASN:H	1:A:110:ASN:ND2	1.78	0.81
1:B:3:LEU:HD22	1:B:4:LYS:H	1.46	0.81
1:A:421:ALA:HB2	1:B:399:ILE:HD11	1.62	0.80
1:C:127:VAL:HA	1:C:236:VAL:HG23	1.64	0.79
1:A:40:VAL:HG21	1:A:76:LEU:HD21	1.62	0.79
1:B:390:VAL:HG12	1:B:395:ILE:HD12	1.64	0.79
1:A:37:THR:HG23	1:A:39:TRP:H	1.47	0.78
1:A:189:ASP:H	1:A:389:GLN:NE2	1.79	0.78
1:A:226:THR:HG23	1:A:228:ALA:H	1.48	0.78
1:C:289:THR:HG23	1:C:291:ARG:HD2	1.65	0.77
1:A:189:ASP:N	1:A:389:GLN:HE22	1.78	0.77
1:C:361:THR:HG22	1:C:370:ARG:H	1.50	0.76
1:C:220:ARG:HB3	1:C:220:ARG:NH1	2.00	0.75
1:B:111:GLY:HA2	1:B:280:ASP:HB2	1.70	0.74
1:A:289:THR:HG22	1:A:291:ARG:H	1.52	0.74
1:C:148:VAL:HG13	1:C:234:GLU:HB2	1.70	0.73
1:A:390:VAL:HG12	1:A:395:ILE:HD12	1.69	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:295:VAL:HG12	1:C:297:LEU:HD12	1.69	0.73
1:C:232:LYS:HZ3	1:C:233:ALA:H	1.37	0.73
1:C:289:THR:HG22	1:C:291:ARG:H	1.54	0.73
1:A:378:VAL:HG21	1:A:440:LEU:HG	1.71	0.72
1:C:84:GLU:HA	1:C:251:GLN:NE2	2.03	0.71
1:A:361:THR:HG23	1:A:362:ARG:HD3	1.73	0.71
1:B:85:VAL:H	1:B:251:GLN:NE2	1.88	0.71
1:A:124:LEU:HD23	1:A:221:VAL:HA	1.73	0.70
1:B:11:GLY:HA3	1:B:35:GLU:HG2	1.72	0.70
1:B:126:GLY:HA3	1:B:221:VAL:CB	2.23	0.69
1:A:148:VAL:HG13	1:A:234:GLU:HB2	1.73	0.69
1:C:232:LYS:NZ	1:C:233:ALA:H	1.90	0.69
1:A:152:VAL:HB	1:A:236:VAL:HG12	1.75	0.69
1:B:132:LEU:C	1:B:134:PRO:HD2	2.14	0.68
1:B:265:ASN:HD21	1:B:269:GLN:NE2	1.91	0.68
1:B:265:ASN:HD21	1:B:269:GLN:HE21	1.40	0.68
1:B:341:THR:HG22	1:B:344:GLU:H	1.59	0.68
1:C:289:THR:CG2	1:C:291:ARG:HD2	2.25	0.67
1:A:210:GLU:OE2	1:A:226:THR:HG21	1.96	0.66
1:A:264:THR:HG23	1:A:268:MET:HA	1.77	0.66
1:A:397:PRO:HG3	1:B:397:PRO:HG3	1.78	0.66
1:A:6:VAL:HG21	1:A:22:VAL:HG21	1.77	0.66
1:A:341:THR:HG22	1:A:344:GLU:H	1.61	0.66
1:B:243:LYS:HG2	1:B:244:PRO:HD2	1.79	0.65
1:A:127:VAL:HG22	1:A:236:VAL:CG2	2.26	0.65
1:B:133:PRO:N	1:B:134:PRO:HD2	2.11	0.65
1:B:243:LYS:HG2	1:B:244:PRO:CD	2.27	0.64
1:C:134:PRO:O	1:C:138:ALA:N	2.30	0.64
1:B:268:MET:HG2	1:B:307:VAL:CG2	2.27	0.64
1:B:148:VAL:HG11	1:B:235:LEU:HB2	1.80	0.64
1:B:420:LEU:HD12	1:B:434:ILE:CD1	2.28	0.63
1:B:131:ASP:OD1	1:B:132:LEU:N	2.32	0.63
1:C:146:TYR:O	1:C:148:VAL:HG23	1.98	0.63
1:C:181:GLU:CD	1:C:181:GLU:H	2.02	0.62
1:B:341:THR:CG2	1:B:344:GLU:H	2.11	0.62
1:B:288:ILE:HD11	1:B:406:LEU:HG	1.82	0.62
1:B:181:GLU:CD	1:B:181:GLU:H	2.02	0.61
1:C:128:PHE:O	1:C:237:ILE:HA	2.00	0.61
1:C:125:LYS:HB2	1:C:220:ARG:HA	1.83	0.61
1:A:284:THR:HB	1:A:304:MET:HE3	1.81	0.61
1:A:38:GLU:HG3	1:A:79:ASN:HD21	1.66	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:215:ILE:HG22	1:C:224:VAL:HG23	1.83	0.61
1:B:268:MET:HG2	1:B:307:VAL:HG22	1.81	0.60
1:B:40:VAL:HG21	1:B:76:LEU:HD21	1.84	0.60
1:A:36:ALA:O	1:A:79:ASN:HA	2.01	0.60
1:A:353:ARG:HB3	1:A:380:ASN:ND2	2.16	0.60
1:B:375:LYS:HB3	1:B:389:GLN:HB2	1.83	0.60
1:C:111:GLY:N	1:C:281:VAL:HG22	2.17	0.60
1:C:148:VAL:HA	1:C:234:GLU:HG2	1.82	0.60
1:B:148:VAL:HG13	1:B:234:GLU:HB2	1.84	0.60
1:C:361:THR:HG23	1:C:362:ARG:HD3	1.83	0.60
1:A:37:THR:HG1	1:A:39:TRP:HE3	1.50	0.59
1:A:288:ILE:HD11	1:A:406:LEU:HG	1.84	0.59
1:A:180:GLY:HA3	1:A:185:ARG:HE	1.67	0.59
1:B:128:PHE:O	1:B:237:ILE:HA	2.02	0.59
1:A:16:MET:CE	1:A:33:VAL:HG11	2.32	0.59
1:A:378:VAL:CG2	1:A:440:LEU:HG	2.32	0.59
1:A:40:VAL:HG23	1:A:78:LEU:HD21	1.84	0.59
1:A:268:MET:HG2	1:A:307:VAL:CG2	2.33	0.58
1:C:166:ALA:O	1:C:170:GLN:HG2	2.03	0.58
1:A:159:ILE:HG13	1:A:239:ALA:HB1	1.84	0.58
1:C:265:ASN:HD21	1:C:269:GLN:HE21	1.51	0.58
1:C:375:LYS:HB3	1:C:389:GLN:HB2	1.86	0.58
1:A:16:MET:HE2	1:A:33:VAL:HG11	1.86	0.57
1:B:278:ALA:HB1	1:B:304:MET:HB3	1.85	0.57
1:A:361:THR:HG22	1:A:370:ARG:H	1.70	0.57
1:C:361:THR:CG2	1:C:370:ARG:H	2.16	0.57
1:B:3:LEU:HD13	1:B:4:LYS:N	2.19	0.57
1:B:420:LEU:HD12	1:B:434:ILE:HD11	1.84	0.57
1:C:16:MET:HE3	1:C:33:VAL:HG11	1.87	0.57
1:C:285:ARG:HH11	1:C:285:ARG:HB3	1.71	0.56
1:A:341:THR:CG2	1:A:344:GLU:H	2.18	0.56
1:C:83:ILE:O	1:C:84:GLU:O	2.22	0.56
1:C:390:VAL:HG12	1:C:395:ILE:HD12	1.87	0.56
1:A:336:GLY:HA3	3:A:454:HOH:O	2.06	0.56
1:B:390:VAL:HB	1:B:395:ILE:HG23	1.88	0.55
1:B:438:ARG:O	1:B:442:GLU:HG3	2.06	0.55
1:C:132:LEU:C	1:C:134:PRO:HD2	2.27	0.55
1:B:341:THR:HG23	1:B:343:MET:H	1.72	0.55
1:B:212:THR:HA	1:B:226:THR:HB	1.87	0.55
1:B:361:THR:HG22	1:B:369:GLY:HA2	1.89	0.55
1:B:142:TYR:O	1:B:145:LYS:HB3	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:268:MET:HE2	1:A:304:MET:HB3	1.89	0.55
1:B:261:ALA:HB3	1:B:283:GLU:HB2	1.87	0.55
1:C:398:ARG:H	1:C:398:ARG:HD2	1.71	0.54
1:C:215:ILE:HG13	1:C:215:ILE:O	2.06	0.54
1:C:242:ILE:HD11	3:C:454:HOH:O	2.06	0.54
1:C:268:MET:HG2	1:C:307:VAL:CG2	2.38	0.54
1:A:421:ALA:CB	1:B:399:ILE:HD11	2.36	0.54
1:B:244:PRO:HG3	1:B:294:TRP:CZ2	2.43	0.54
1:C:16:MET:CE	1:C:33:VAL:HG11	2.37	0.54
1:C:232:LYS:HG3	1:C:233:ALA:N	2.22	0.54
1:A:3:LEU:HD11	1:A:5:LYS:NZ	2.22	0.54
1:A:334:GLU:HB2	1:A:395:ILE:CG1	2.38	0.54
1:A:408:ALA:HB2	1:B:410:PHE:CE1	2.43	0.54
1:B:125:LYS:HB2	1:B:219:GLU:O	2.07	0.54
1:A:264:THR:O	1:A:292:ARG:NH1	2.41	0.53
1:C:16:MET:HE1	1:C:33:VAL:HG21	1.90	0.53
1:A:302:ASN:HB3	1:B:430:TRP:CZ2	2.44	0.53
1:C:289:THR:HG22	1:C:291:ARG:HB2	1.91	0.53
1:B:37:THR:HG23	1:B:38:GLU:N	2.22	0.53
1:A:110:ASN:HD22	1:A:110:ASN:N	1.99	0.53
1:A:119:ILE:O	1:A:122:VAL:HG23	2.08	0.53
1:C:365:TYR:CD1	1:C:366:TYR:N	2.77	0.53
1:A:93:ARG:HH11	1:A:93:ARG:CB	2.12	0.53
1:A:353:ARG:HB3	1:A:380:ASN:HD22	1.73	0.52
1:B:364:HIS:O	1:B:364:HIS:ND1	2.43	0.52
1:A:373:TRP:HB2	1:A:391:VAL:HG13	1.90	0.52
1:B:4:LYS:HB2	1:B:29:TRP:CD2	2.44	0.52
1:C:341:THR:CG2	1:C:344:GLU:H	2.23	0.52
1:C:398:ARG:HD2	1:C:398:ARG:N	2.25	0.52
1:A:295:VAL:HG12	1:A:297:LEU:HD13	1.92	0.52
1:A:361:THR:O	1:A:362:ARG:HD2	2.10	0.52
1:A:81:GLU:HB3	1:A:93:ARG:HH12	1.75	0.52
1:A:213:MET:HE1	1:A:227:ASP:HA	1.92	0.52
1:C:84:GLU:HA	1:C:251:GLN:HE22	1.75	0.52
1:B:178:VAL:HG11	1:B:183:VAL:HG12	1.91	0.52
1:A:334:GLU:HB2	1:A:395:ILE:HG13	1.93	0.51
1:C:85:VAL:H	1:C:251:GLN:HE21	1.56	0.51
1:A:35:GLU:CD	1:A:37:THR:HG22	2.30	0.51
1:B:199:LYS:HG2	1:B:333:VAL:CG1	2.40	0.51
1:B:91:ARG:HD3	1:B:98:GLU:OE2	2.10	0.51
1:C:189:ASP:N	1:C:389:GLN:HE22	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:ARG:HD3	1:B:98:GLU:CD	2.31	0.51
1:B:133:PRO:O	1:B:136:ALA:HB3	2.10	0.51
1:C:374:LEU:HG	1:C:390:VAL:HG13	1.93	0.51
1:B:189:ASP:N	1:B:389:GLN:HE22	1.98	0.51
1:B:84:GLU:CA	1:B:251:GLN:HE22	2.18	0.51
1:C:264:THR:HG22	1:C:283:GLU:O	2.11	0.51
1:A:398:ARG:HD2	1:A:398:ARG:N	2.25	0.51
1:A:264:THR:CG2	1:A:268:MET:HA	2.40	0.50
1:C:142:TYR:OH	1:C:220:ARG:HB2	2.11	0.50
1:C:361:THR:HG22	1:C:369:GLY:HA2	1.92	0.50
1:A:23:LYS:HE2	1:A:73:GLY:HA3	1.94	0.50
1:C:207:ARG:HH11	1:C:207:ARG:HA	1.77	0.50
1:B:211:ILE:N	1:B:211:ILE:HD12	2.26	0.50
1:B:3:LEU:HD22	1:B:4:LYS:N	2.22	0.50
1:A:127:VAL:HG13	1:A:238:LEU:HD13	1.92	0.50
1:C:154:ILE:CD1	1:C:212:THR:HG21	2.42	0.50
1:A:202:LYS:HD2	3:A:567:HOH:O	2.11	0.50
1:C:11:GLY:HA3	1:C:35:GLU:HG2	1.94	0.50
1:C:20:SER:HA	1:C:74:ILE:HD11	1.94	0.50
1:C:364:HIS:HA	3:C:467:HOH:O	2.11	0.49
1:A:189:ASP:OD1	1:A:343:MET:HG3	2.12	0.49
1:C:289:THR:CG2	1:C:291:ARG:HB2	2.43	0.49
1:C:285:ARG:HH11	1:C:285:ARG:CB	2.24	0.49
1:C:383:ASN:HD22	1:C:412:THR:HG22	1.77	0.49
1:A:420:LEU:HD12	1:A:434:ILE:HD13	1.95	0.49
1:A:291:ARG:CZ	1:C:291:ARG:HH11	2.25	0.49
1:C:84:GLU:HG2	3:C:452:HOH:O	2.13	0.49
1:C:323:LEU:HD21	1:C:406:LEU:HB3	1.95	0.49
1:A:300:ALA:O	1:A:304:MET:HG3	2.11	0.49
1:C:334:GLU:HB2	1:C:395:ILE:CG1	2.42	0.49
1:C:3:LEU:HD13	1:C:3:LEU:C	2.32	0.49
1:C:242:ILE:HG23	1:C:294:TRP:CH2	2.48	0.49
1:A:261:ALA:HB1	1:A:281:VAL:O	2.13	0.49
1:B:127:VAL:CA	1:B:236:VAL:HG23	2.39	0.49
1:C:4:LYS:HB2	1:C:29:TRP:CD2	2.48	0.48
1:A:390:VAL:HG12	1:A:395:ILE:CD1	2.42	0.48
1:C:412:THR:HG21	1:C:441:LYS:HA	1.94	0.48
1:B:341:THR:HG23	1:B:343:MET:N	2.28	0.48
1:C:364:HIS:O	1:C:365:TYR:HB3	2.12	0.48
1:B:257:GLY:HA3	1:B:283:GLU:OE1	2.12	0.48
1:B:398:ARG:H	1:B:398:ARG:HD2	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:431:ASP:O	1:C:435:VAL:HG23	2.13	0.48
1:A:289:THR:CG2	1:A:291:ARG:HB2	2.43	0.48
1:C:372:ILE:HD12	1:C:393:SER:O	2.14	0.48
1:B:334:GLU:HB2	1:B:395:ILE:CG1	2.43	0.48
1:B:427:ALA:HB1	1:B:428:PRO:HD2	1.96	0.48
1:C:152:VAL:HG11	1:C:224:VAL:HG11	1.96	0.48
1:C:341:THR:HG23	1:C:344:GLU:H	1.78	0.48
1:A:132:LEU:C	1:A:134:PRO:HD2	2.35	0.48
1:B:16:MET:HE1	1:B:76:LEU:HD13	1.95	0.48
1:C:242:ILE:HG22	3:C:485:HOH:O	2.13	0.48
1:C:334:GLU:HB2	1:C:395:ILE:HG13	1.96	0.48
1:A:153:ILE:HG12	1:A:237:ILE:HB	1.96	0.47
1:C:179:ARG:O	1:C:209:GLN:HA	2.14	0.47
1:B:211:ILE:HG22	1:B:212:THR:O	2.13	0.47
1:B:224:VAL:CG1	1:B:231:TYR:HB2	2.44	0.47
1:A:396:LEU:HB3	1:A:397:PRO:HD3	1.96	0.47
1:B:193:THR:HG22	1:B:197:GLU:HG3	1.97	0.47
1:A:18:ALA:O	1:A:22:VAL:HG12	2.15	0.47
1:A:410:PHE:CE1	1:B:408:ALA:HB2	2.50	0.47
1:B:378:VAL:HG21	1:B:440:LEU:HG	1.96	0.47
1:B:438:ARG:HD2	1:B:442:GLU:OE2	2.13	0.47
1:C:214:LYS:O	1:C:225:VAL:HG12	2.15	0.47
1:B:361:THR:HG22	1:B:370:ARG:H	1.79	0.47
1:C:220:ARG:HH11	1:C:220:ARG:CB	2.17	0.47
1:C:427:ALA:HB1	1:C:428:PRO:HD2	1.97	0.47
1:C:361:THR:HG22	1:C:370:ARG:N	2.27	0.47
1:A:37:THR:HG23	1:A:39:TRP:N	2.24	0.46
1:A:267:LYS:HD3	1:A:316:GLU:OE2	2.15	0.46
1:A:327:VAL:HG13	1:B:423:ALA:HB2	1.97	0.46
1:B:150:ASN:HD22	1:B:232:LYS:HB3	1.80	0.46
1:B:232:LYS:HA	1:B:232:LYS:HE2	1.97	0.46
1:B:384:ARG:HE	1:B:384:ARG:HB2	1.55	0.46
1:B:138:ALA:O	1:B:139:ILE:C	2.52	0.46
1:C:341:THR:HG22	1:C:344:GLU:HB2	1.97	0.46
1:B:289:THR:HB	1:B:291:ARG:HG2	1.96	0.46
1:B:334:GLU:HB2	1:B:395:ILE:HG12	1.98	0.46
1:A:83:ILE:HG22	1:A:91:ARG:O	2.15	0.46
1:A:398:ARG:HD2	1:A:398:ARG:H	1.80	0.46
1:B:423:ALA:HB3	1:B:426:PHE:HD2	1.81	0.46
1:A:181:GLU:CD	1:A:181:GLU:H	2.19	0.46
1:A:303:LYS:O	1:A:307:VAL:HG13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:VAL:HG22	1:B:276:TYR:HB2	1.97	0.46
1:C:235:LEU:HD13	1:C:236:VAL:N	2.31	0.46
1:C:244:PRO:HB2	1:C:246:ILE:HG23	1.97	0.46
1:B:80:ALA:HA	1:B:94:GLU:HB2	1.97	0.46
1:B:189:ASP:H	1:B:389:GLN:NE2	1.99	0.46
1:C:242:ILE:HG23	1:C:294:TRP:HH2	1.80	0.46
1:C:399:ILE:HG13	1:C:400:ASP:N	2.31	0.46
1:A:361:THR:HG22	1:A:369:GLY:HA2	1.97	0.45
1:A:421:ALA:HB2	1:B:399:ILE:CD1	2.41	0.45
1:A:341:THR:HG23	1:A:343:MET:H	1.81	0.45
1:B:143:MET:HA	1:B:148:VAL:CG2	2.47	0.45
1:A:4:LYS:HG3	1:A:29:TRP:CZ2	2.51	0.45
1:A:289:THR:HG22	1:A:291:ARG:HB2	1.98	0.45
1:A:361:THR:CG2	1:A:370:ARG:H	2.29	0.45
1:C:174:VAL:O	1:C:204:VAL:HG13	2.17	0.45
1:A:119:ILE:HD11	1:A:215:ILE:HD11	1.97	0.45
1:C:265:ASN:HD21	1:C:269:GLN:NE2	2.14	0.45
1:B:361:THR:CG2	1:B:370:ARG:H	2.30	0.45
1:C:213:MET:HE1	1:C:227:ASP:HA	1.98	0.45
1:A:284:THR:HB	1:A:304:MET:CE	2.45	0.45
1:C:37:THR:HG23	1:C:38:GLU:N	2.32	0.45
1:B:431:ASP:HB3	1:B:434:ILE:HD13	1.98	0.44
1:C:128:PHE:CE2	1:C:139:ILE:HA	2.52	0.44
1:A:112:ALA:HB3	1:A:242:ILE:HD11	1.98	0.44
1:A:373:TRP:HB2	1:A:391:VAL:CG1	2.46	0.44
1:C:143:MET:C	1:C:145:LYS:H	2.21	0.44
1:B:390:VAL:CG1	1:B:395:ILE:HD12	2.42	0.44
1:A:3:LEU:C	1:A:3:LEU:HD13	2.38	0.44
1:B:85:VAL:H	1:B:251:GLN:HE21	1.63	0.44
1:A:82:VAL:HG13	1:A:90:VAL:HG22	2.00	0.44
1:C:207:ARG:HA	1:C:207:ARG:NH1	2.32	0.44
1:A:296:PRO:C	1:A:297:LEU:HD12	2.38	0.44
1:B:143:MET:HA	1:B:148:VAL:HG23	2.00	0.44
1:B:378:VAL:CG2	1:B:440:LEU:HG	2.47	0.44
1:A:35:GLU:OE1	1:A:36:ALA:N	2.52	0.43
1:A:122:VAL:HG13	1:A:238:LEU:HD21	2.00	0.43
1:B:287:VAL:HG21	1:B:407:MET:CE	2.49	0.43
1:A:399:ILE:HG13	1:A:400:ASP:N	2.33	0.43
1:A:361:THR:CG2	1:A:362:ARG:HD3	2.46	0.43
1:C:131:ASP:OD1	1:C:132:LEU:N	2.51	0.43
1:A:365:TYR:CD1	1:A:366:TYR:N	2.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:THR:CG2	1:B:39:TRP:H	2.14	0.43
1:A:289:THR:HG22	1:A:291:ARG:N	2.28	0.43
1:B:267:LYS:O	1:B:268:MET:HB2	2.19	0.43
1:A:423:ALA:HB3	1:A:426:PHE:HD2	1.84	0.43
1:A:438:ARG:HD2	1:A:442:GLU:CD	2.39	0.43
1:B:303:LYS:O	1:B:307:VAL:HG13	2.19	0.43
1:A:8:ILE:HD12	1:A:19:ALA:HB2	2.00	0.43
1:B:35:GLU:C	1:B:37:THR:H	2.21	0.43
1:B:132:LEU:CB	1:B:134:PRO:HD2	2.49	0.43
1:A:244:PRO:HG3	1:A:294:TRP:CZ2	2.53	0.42
1:B:39:TRP:CH2	1:B:134:PRO:HB3	2.54	0.42
1:C:365:TYR:CG	1:C:366:TYR:N	2.87	0.42
1:B:111:GLY:CA	1:B:280:ASP:HB2	2.44	0.42
1:A:81:GLU:HB3	1:A:93:ARG:NH1	2.32	0.42
1:B:170:GLN:HG2	3:B:472:HOH:O	2.19	0.42
1:C:424:PRO:HB2	1:C:425:PRO:HD3	2.01	0.42
1:A:81:GLU:O	1:A:92:VAL:HG23	2.17	0.42
1:A:427:ALA:HB1	1:A:428:PRO:HD2	2.02	0.42
1:C:39:TRP:CH2	1:C:134:PRO:HD3	2.54	0.42
1:A:273:GLU:HG2	1:A:274:ASN:ND2	2.35	0.42
1:A:430:TRP:CZ2	1:B:302:ASN:HB3	2.54	0.42
1:A:429:VAL:HG11	1:B:302:ASN:ND2	2.35	0.42
1:B:4:LYS:HG3	1:B:29:TRP:CZ2	2.55	0.42
1:B:398:ARG:HD2	1:B:398:ARG:N	2.35	0.42
1:C:95:ASN:C	1:C:97:GLY:H	2.23	0.42
1:B:127:VAL:HG22	1:B:236:VAL:HG21	2.01	0.42
1:C:412:THR:CG2	1:C:441:LYS:HA	2.50	0.42
1:A:401:THR:OG1	1:B:400:ASP:HB3	2.19	0.42
1:B:83:ILE:HG13	1:B:83:ILE:O	2.20	0.42
1:C:150:ASN:CG	1:C:232:LYS:HG2	2.40	0.42
1:B:307:VAL:HA	1:B:317:LEU:HD23	2.02	0.41
1:C:112:ALA:HB1	1:C:242:ILE:HG12	2.01	0.41
1:A:84:GLU:OE1	1:A:91:ARG:HD2	2.20	0.41
1:C:179:ARG:HA	1:C:210:GLU:H	1.85	0.41
1:C:390:VAL:HB	1:C:395:ILE:HG23	2.01	0.41
1:A:76:LEU:HD12	1:A:77:HIS:N	2.35	0.41
1:A:188:PHE:HA	1:A:389:GLN:NE2	2.35	0.41
1:A:408:ALA:HB2	1:B:410:PHE:HE1	1.82	0.41
1:B:127:VAL:HG22	1:B:236:VAL:CG2	2.49	0.41
1:B:150:ASN:ND2	1:B:232:LYS:HB3	2.35	0.41
1:C:38:GLU:HB3	1:C:79:ASN:HD21	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:154:ILE:HD13	1:C:212:THR:HG21	2.01	0.41
1:B:133:PRO:N	1:B:134:PRO:CD	2.81	0.41
1:C:287:VAL:HG21	1:C:407:MET:CE	2.50	0.41
1:A:40:VAL:HG21	1:A:76:LEU:CD2	2.43	0.41
1:B:35:GLU:OE1	1:B:37:THR:N	2.53	0.41
1:C:383:ASN:HD22	1:C:412:THR:CG2	2.34	0.41
1:A:152:VAL:HG11	1:A:224:VAL:HG11	2.02	0.41
1:A:159:ILE:CG1	1:A:239:ALA:HB1	2.50	0.41
1:A:161:ILE:HD13	1:A:196:LEU:HD21	2.03	0.41
1:A:264:THR:CG2	1:A:265:ASN:N	2.83	0.41
1:C:330:PHE:O	1:C:333:VAL:HG12	2.20	0.41
1:A:22:VAL:HG13	1:A:23:LYS:N	2.36	0.41
1:B:215:ILE:CG2	1:B:224:VAL:HA	2.42	0.41
1:C:114:PRO:HG3	1:C:131:ASP:HB2	2.03	0.41
1:C:133:PRO:O	1:C:136:ALA:HB3	2.21	0.41
1:A:224:VAL:HG12	1:A:231:TYR:HB2	2.03	0.41
1:A:341:THR:HG23	1:A:343:MET:N	2.36	0.41
1:C:8:ILE:HG12	1:C:107:VAL:HB	2.02	0.41
1:C:435:VAL:O	1:C:439:VAL:HG23	2.20	0.41
1:A:151:VAL:HG11	1:A:167:PHE:HB3	2.03	0.40
1:A:410:PHE:HE1	1:B:408:ALA:HB2	1.85	0.40
1:A:357:ILE:HD13	1:A:435:VAL:HG12	2.02	0.40
1:B:124:LEU:C	1:B:126:GLY:H	2.24	0.40
1:C:35:GLU:OE1	1:C:37:THR:N	2.40	0.40
1:B:3:LEU:O	1:B:4:LYS:HG2	2.21	0.40
1:B:112:ALA:HB1	1:B:242:ILE:HG12	2.03	0.40
1:A:246:ILE:O	1:A:250:LYS:HG3	2.21	0.40
1:B:26:LYS:HB3	1:B:29:TRP:CG	2.57	0.40
1:B:289:THR:HG22	1:B:291:ARG:NH2	2.36	0.40
1:C:84:GLU:CA	1:C:251:GLN:HE22	2.35	0.40
1:A:101:TYR:CD2	1:A:101:TYR:N	2.89	0.40
1:C:154:ILE:HD11	1:C:212:THR:HG21	2.04	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	414/449 (92%)	384 (93%)	26 (6%)	4 (1%)	15 27
1	B	414/449 (92%)	375 (91%)	29 (7%)	10 (2%)	6 9
1	C	414/449 (92%)	373 (90%)	32 (8%)	9 (2%)	6 11
All	All	1242/1347 (92%)	1132 (91%)	87 (7%)	23 (2%)	8 14

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	VAL
1	A	145	LYS
1	B	138	ALA
1	C	84	GLU
1	C	365	TYR
1	A	218	GLU
1	B	121	GLY
1	B	145	LYS
1	B	228	ALA
1	C	67	VAL
1	C	227	ASP
1	B	139	ILE
1	B	214	LYS
1	B	227	ASP
1	C	144	GLU
1	C	228	ALA
1	B	125	LYS
1	C	87	THR
1	B	10	GLY
1	A	10	GLY
1	C	10	GLY
1	B	67	VAL
1	C	296	PRO

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	314/364 (86%)	289 (92%)	25 (8%)	12 21
1	B	315/364 (86%)	276 (88%)	39 (12%)	4 7
1	C	314/364 (86%)	281 (90%)	33 (10%)	7 11
All	All	943/1092 (86%)	846 (90%)	97 (10%)	7 12

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	GLU
1	A	86	ASP
1	A	93	ARG
1	A	98	GLU
1	A	110	ASN
1	A	198	GLU
1	A	205	ASN
1	A	224	VAL
1	A	234	GLU
1	A	252	LEU
1	A	256	ILE
1	A	281	VAL
1	A	285	ARG
1	A	316	GLU
1	A	323	LEU
1	A	341	THR
1	A	361	THR
1	A	364	HIS
1	A	394	ASP
1	A	398	ARG
1	A	399	ILE
1	A	406	LEU
1	A	413	LYS
1	A	438	ARG
1	A	440	LEU
1	B	3	LEU
1	B	28	GLU
1	B	35	GLU
1	B	37	THR
1	B	86	ASP
1	B	98	GLU

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Mol	Chain	Res	Type
1	B	146	TYR
1	B	159	ILE
1	B	170	GLN
1	B	173	ASN
1	B	178	VAL
1	B	205	ASN
1	B	209	GLN
1	B	215	ILE
1	B	226	THR
1	B	232	LYS
1	B	235	LEU
1	B	236	VAL
1	B	238	LEU
1	B	252	LEU
1	B	264	THR
1	B	281	VAL
1	B	285	ARG
1	B	289	THR
1	B	291	ARG
1	B	297	LEU
1	B	307	VAL
1	B	315	LYS
1	B	323	LEU
1	B	333	VAL
1	B	361	THR
1	B	364	HIS
1	B	383	ASN
1	B	394	ASP
1	B	398	ARG
1	B	399	ILE
1	B	406	LEU
1	B	438	ARG
1	B	440	LEU
1	C	28	GLU
1	C	35	GLU
1	C	37	THR
1	C	83	ILE
1	C	84	GLU
1	C	89	TYR
1	C	91	ARG
1	C	92	VAL
1	C	98	GLU

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Mol	Chain	Res	Type
1	C	142	TYR
1	C	178	VAL
1	C	181	GLU
1	C	207	ARG
1	C	216	GLU
1	C	220	ARG
1	C	226	THR
1	C	232	LYS
1	C	234	GLU
1	C	238	LEU
1	C	252	LEU
1	C	281	VAL
1	C	285	ARG
1	C	291	ARG
1	C	316	GLU
1	C	347	LYS
1	C	362	ARG
1	C	364	HIS
1	C	394	ASP
1	C	398	ARG
1	C	399	ILE
1	C	406	LEU
1	C	434	ILE
1	C	438	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	95	ASN
1	A	110	ASN
1	A	205	ASN
1	A	269	GLN
1	A	274	ASN
1	A	302	ASN
1	A	380	ASN
1	A	389	GLN
1	B	95	ASN
1	B	115	GLN
1	B	150	ASN
1	B	205	ASN
1	B	209	GLN

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Mol	Chain	Res	Type
1	B	251	GLN
1	B	269	GLN
1	B	302	ASN
1	B	380	ASN
1	B	389	GLN
1	C	79	ASN
1	C	115	GLN
1	C	205	ASN
1	C	251	GLN
1	C	269	GLN
1	C	302	ASN
1	C	380	ASN
1	C	383	ASN
1	C	389	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is 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	GOL	A	968	-	5,5,5	0.51	0	5,5,5	0.43	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	GOL	A	968	-	-	0/4/4/4	-

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	418/449 (93%)	0.09	8 (1%)	66	75	28, 51, 70, 77	0
1	B	418/449 (93%)	0.11	10 (2%)	59	68	30, 50, 71, 83	0
1	C	418/449 (93%)	0.20	11 (2%)	56	65	28, 51, 73, 82	0
All	All	1254/1347 (93%)	0.13	29 (2%)	60	69	28, 51, 72, 83	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	134	PRO	6.1
1	B	134	PRO	4.8
1	C	136	ALA	4.6
1	A	134	PRO	4.2
1	A	93	ARG	4.1
1	A	91	ARG	3.7
1	C	215	ILE	3.3
1	B	235	LEU	3.3
1	C	140	ARG	3.2
1	C	144	GLU	3.0
1	A	80	ALA	2.9
1	C	217	GLY	2.9
1	B	364	HIS	2.8
1	C	143	MET	2.7
1	C	238	LEU	2.6
1	A	215	ILE	2.5
1	C	224	VAL	2.5
1	C	145	LYS	2.4
1	B	219	GLU	2.4
1	A	133	PRO	2.4
1	B	136	ALA	2.3
1	B	2	SER	2.3
1	B	123	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	215	ILE	2.2
1	A	39	TRP	2.2
1	C	212	THR	2.2
1	A	151	VAL	2.1
1	B	120	GLU	2.1
1	B	126	GLY	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 monosaccharides 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	GOL	A	968	6/6	0.84	0.23	58,59,60,60	0

6.5 Other polymers (i)

There are no such residues in this entry.