



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 7, 2023 – 12:21 AM EDT

PDB ID : 4G5E  
Title : 2,4,6-Trichlorophenol 4-monoxygenase  
Authors : Hayes, R.P.; Webb, B.N.; Subramanian, A.K.; Nissen, M.; Popchock, A.; Xun, L.; Kang, C.  
Deposited on : 2012-07-17  
Resolution : 2.50 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](mailto: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>.

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.35  
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) : 2.35

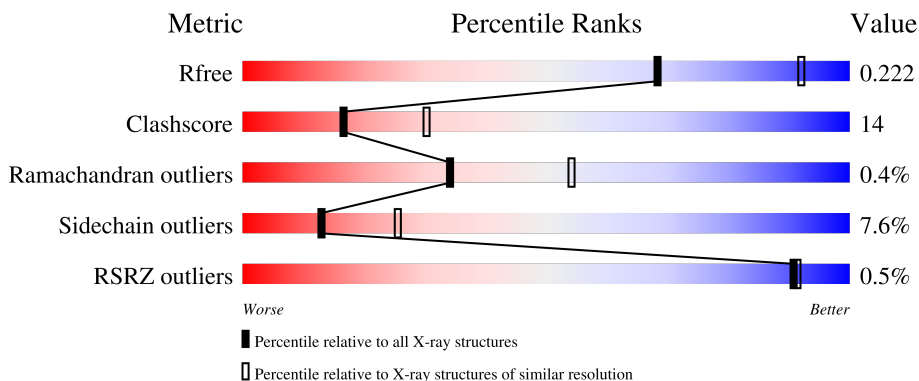
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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  $\geq 3$ , 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	517	 66% 21% 9%
1	B	517	 62% 27% 9%
1	C	517	 67% 22% 9%
1	D	517	 60% 27% 9%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15418 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 2,4,6-Trichlorophenol 4-monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	471	3777	2412	670	680	15	0	0	0
1	B	471	3772	2409	670	678	15	0	0	1
1	C	471	3777	2412	670	680	15	0	0	0
1	D	471	3777	2412	670	680	15	0	0	0

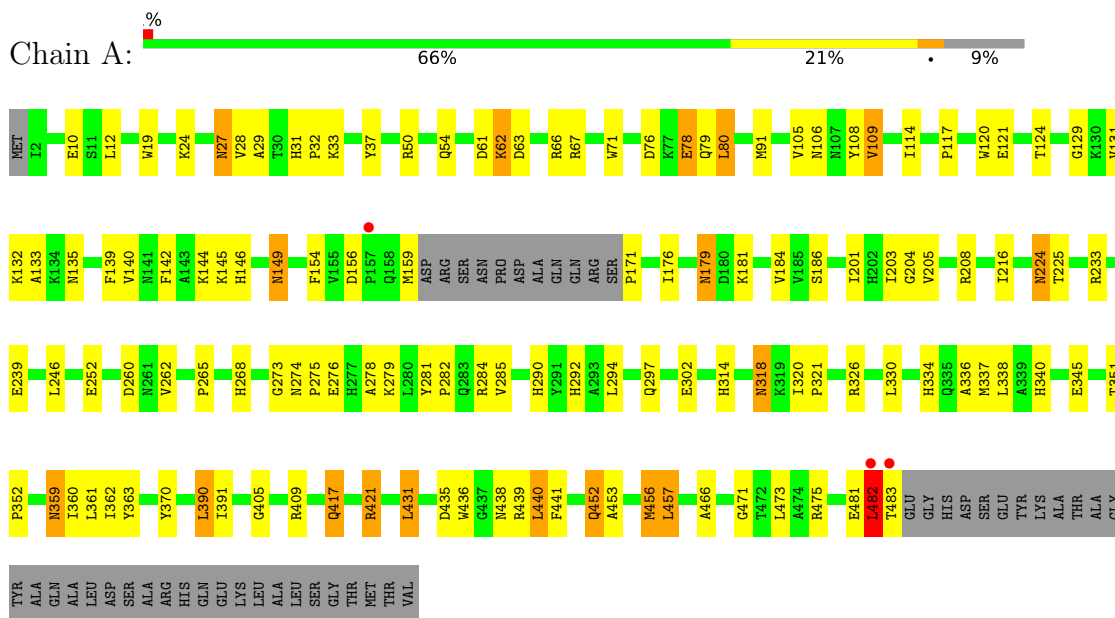
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	92	92	92	0	0
2	B	68	68	68	0	0
2	C	112	112	112	0	0
2	D	43	43	43	0	0

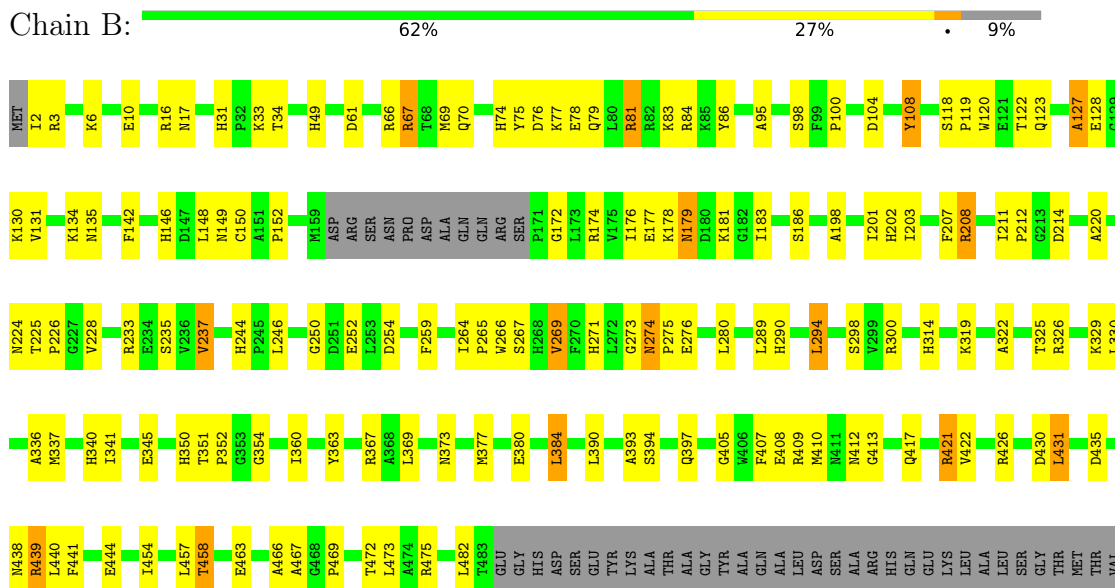
### 3 Residue-property plots

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: 2,4,6-Trichlorophenol 4-monooxygenase

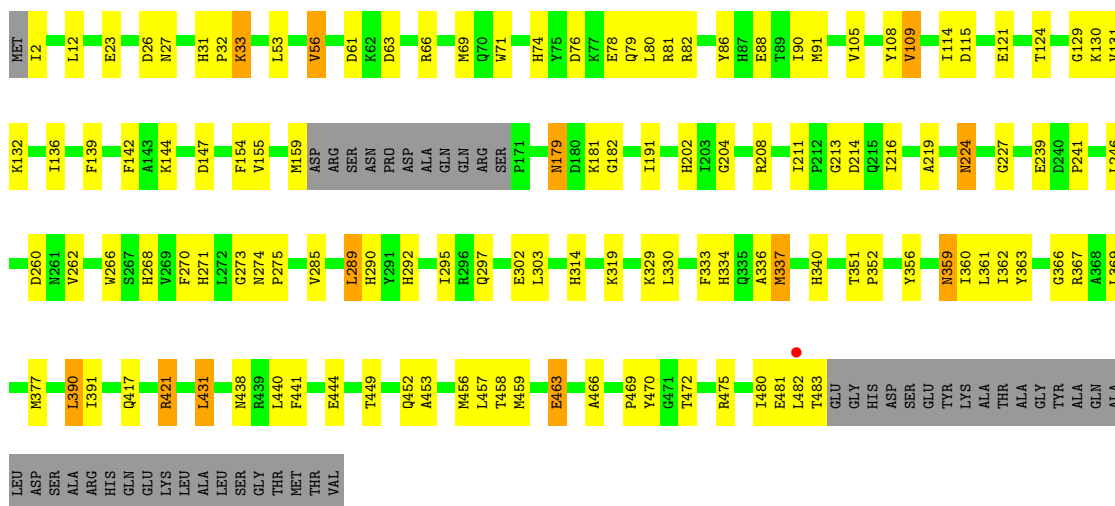


- Molecule 1: 2,4,6-Trichlorophenol 4-monooxygenase



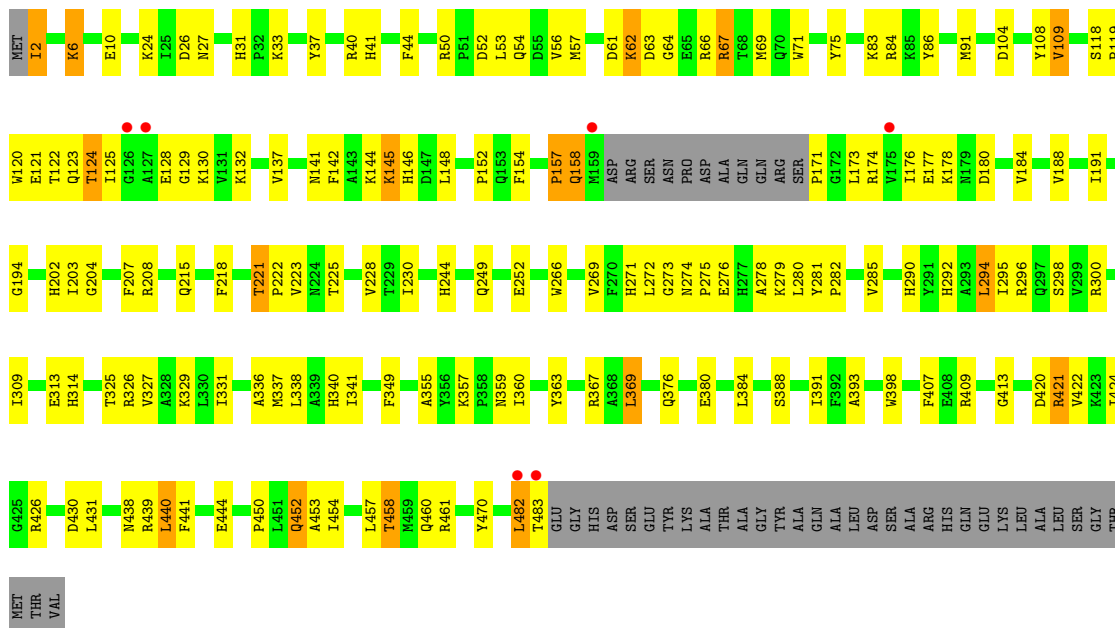
- Molecule 1: 2,4,6-Trichlorophenol 4-monooxygenase

Chain C:  67% 22% 9%



- Molecule 1: 2,4,6-Trichlorophenol 4-monooxygenase

Chain D:  60% 27% 9%



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	107.07Å 180.05Å 110.20Å 90.00° 97.87° 90.00°	Depositor
Resolution (Å)	49.50 – 2.50 49.50 – 2.04	Depositor EDS
% Data completeness (in resolution range)	94.8 (49.50-2.50) 85.6 (49.50-2.04)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.13 (at 2.05Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, $R_{free}$	0.174 , 0.224 0.171 , 0.222	Depositor DCC
$R_{free}$ test set	2000 reflections (1.58%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.1	Xtrriage
Anisotropy	0.185	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 42.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	15418	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

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	0/3877	0.57	0/5259
1	B	0.41	0/3872	0.55	0/5252
1	C	0.45	0/3877	0.59	0/5259
1	D	0.40	0/3877	0.53	0/5259
All	All	0.42	0/15503	0.56	0/21029

There are no bond length outliers.

There are no bond angle outliers.

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	3777	0	3695	94	0
1	B	3772	0	3689	114	0
1	C	3777	0	3695	91	0
1	D	3777	0	3695	125	0
2	A	92	0	0	6	0
2	B	68	0	0	4	0
2	C	112	0	0	5	0
2	D	43	0	0	5	0
All	All	15418	0	14774	424	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 14.

All (424) 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:C:213:GLY:HA2	1:C:216:ILE:HD12	1.41	1.02
1:B:369:LEU:O	1:B:369:LEU:HD23	1.64	0.96
1:C:121:GLU:O	1:C:124:THR:HG22	1.71	0.90
1:A:179:ASN:HD22	1:A:181:LYS:H	1.20	0.88
1:C:390:LEU:HG	1:C:391:ILE:HG13	1.56	0.87
1:D:470:TYR:O	2:D:604:HOH:O	1.92	0.86
1:D:482:LEU:H	1:D:482:LEU:HD23	1.41	0.85
1:B:469:PRO:O	1:B:472:THR:HG23	1.76	0.85
1:C:179:ASN:HD22	1:C:181:LYS:H	1.23	0.83
1:A:121:GLU:O	1:A:124:THR:HG22	1.77	0.83
1:D:314:HIS:CD2	1:D:421:ARG:HD3	2.12	0.83
1:B:179:ASN:HD22	1:B:181:LYS:H	1.28	0.81
1:B:211:ILE:HD12	1:B:212:PRO:HD2	1.62	0.81
1:B:482:LEU:H	1:B:482:LEU:HD23	1.44	0.80
1:A:314:HIS:CD2	1:A:421:ARG:HD3	2.17	0.80
1:D:482:LEU:HD23	1:D:482:LEU:N	1.97	0.79
1:B:330:LEU:HD23	1:B:377:MET:HE2	1.64	0.78
1:A:106:ASN:HD22	1:A:149:ASN:HD21	1.30	0.78
1:C:483:THR:O	2:C:654:HOH:O	2.02	0.78
1:D:109:VAL:HG13	1:D:285:VAL:CG1	2.14	0.78
1:B:314:HIS:CD2	1:B:421:ARG:HD3	2.18	0.77
1:A:336:ALA:O	1:A:340:HIS:HD2	1.68	0.77
1:A:314:HIS:HD2	1:A:421:ARG:HD3	1.49	0.76
1:C:314:HIS:HD2	1:C:421:ARG:HH11	1.33	0.76
1:C:314:HIS:CD2	1:C:421:ARG:HD3	2.20	0.76
1:D:454:ILE:O	1:D:458:THR:HG22	1.87	0.75
1:B:273:GLY:O	1:B:275:PRO:HD3	1.87	0.74
1:A:171:PRO:N	2:A:687:HOH:O	2.22	0.73
1:B:294:LEU:HD13	1:B:341:ILE:HG13	1.70	0.73
1:A:337:MET:HE2	1:A:370:TYR:HB2	1.71	0.73
1:A:314:HIS:HD2	1:A:421:ARG:HH11	1.37	0.72
1:D:314:HIS:HD2	1:D:421:ARG:HH11	1.37	0.72
1:C:481:GLU:HG3	1:C:482:LEU:H	1.55	0.71
1:A:62:LYS:HG2	1:A:63:ASP:N	2.05	0.71
1:C:31:HIS:HD2	1:C:33:LYS:H	1.38	0.71
1:C:336:ALA:O	1:C:340:HIS:HD2	1.75	0.70
1:A:135:ASN:OD1	2:A:601:HOH:O	2.10	0.70
1:A:276:GLU:O	1:A:279:LYS:HG2	1.92	0.70
1:D:53:LEU:O	1:D:54:GLN:HB2	1.92	0.69
1:D:223:VAL:HG13	1:D:230:ILE:HD11	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:ALA:O	1:B:340:HIS:HD2	1.74	0.69
1:D:109:VAL:HG13	1:D:285:VAL:HG13	1.74	0.69
1:B:298:SER:HB3	1:B:337:MET:HB3	1.75	0.69
1:B:360:ILE:HA	1:B:363:TYR:CE2	2.27	0.69
1:A:109:VAL:HG22	1:A:285:VAL:O	1.93	0.69
1:D:349:PHE:CZ	1:D:357:LYS:HD2	2.28	0.68
1:A:439:ARG:HH21	1:A:440:LEU:CD1	2.06	0.68
1:C:481:GLU:HG3	1:C:482:LEU:N	2.09	0.68
1:B:244:HIS:HD2	1:B:430:ASP:OD2	1.76	0.68
1:D:26:ASP:OD1	1:D:27:ASN:N	2.22	0.68
1:D:157:PRO:HG2	1:D:171:PRO:HD2	1.76	0.67
1:A:31:HIS:HD2	1:A:33:LYS:H	1.41	0.67
1:B:274:ASN:C	1:B:274:ASN:HD22	1.98	0.67
1:C:129:GLY:O	1:C:132:LYS:HE2	1.94	0.67
1:D:244:HIS:HD2	1:D:430:ASP:OD2	1.79	0.66
1:B:472:THR:HG22	1:B:475:ARG:NH2	2.10	0.66
1:A:390:LEU:HG	1:A:391:ILE:HG12	1.77	0.66
1:B:369:LEU:HD23	1:B:369:LEU:C	2.16	0.65
1:D:176:ILE:HD11	1:D:184:VAL:CG1	2.27	0.65
1:A:179:ASN:ND2	1:A:181:LYS:H	1.92	0.65
1:A:481:GLU:O	1:A:482:LEU:HB2	1.97	0.65
1:B:75:TYR:H	1:B:79:GLN:NE2	1.94	0.65
1:A:318:ASN:H	1:A:318:ASN:HD22	1.43	0.65
1:C:480:ILE:HG22	1:C:481:GLU:HG2	1.78	0.65
1:D:337:MET:CE	1:D:369:LEU:HD12	2.26	0.65
1:D:109:VAL:HG22	1:D:285:VAL:O	1.97	0.65
1:D:121:GLU:OE1	1:D:132:LYS:HA	1.97	0.65
1:D:146:HIS:HB2	1:D:148:LEU:HD13	1.79	0.64
1:B:482:LEU:H	1:B:482:LEU:CD2	2.11	0.64
1:C:390:LEU:HD23	1:C:390:LEU:H	1.63	0.64
1:D:314:HIS:HD2	1:D:421:ARG:HD3	1.57	0.64
1:B:454:ILE:O	1:B:458:THR:HG23	1.97	0.64
1:A:129:GLY:O	1:A:132:LYS:NZ	2.31	0.63
1:A:80:LEU:HD13	1:A:345:GLU:HG2	1.79	0.63
1:D:266:TRP:HA	1:D:269:VAL:HG12	1.79	0.63
1:A:246:LEU:HD12	1:A:431:LEU:HD13	1.80	0.62
1:B:237:VAL:HG13	1:B:250:GLY:HA2	1.80	0.62
1:C:273:GLY:O	1:C:275:PRO:HD3	2.00	0.62
1:B:274:ASN:ND2	1:B:276:GLU:H	1.98	0.61
1:C:314:HIS:HE1	1:C:417:GLN:O	1.82	0.61
1:A:359:ASN:HD22	1:A:362:ILE:H	1.46	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:179:ASN:HD22	1:C:181:LYS:N	1.97	0.61
1:D:118:SER:O	1:D:122:THR:HG23	2.00	0.61
1:B:176:ILE:HD11	1:B:186:SER:HB2	1.81	0.61
1:B:314:HIS:HD2	1:B:421:ARG:HH11	1.48	0.61
1:B:467:ALA:HB1	1:B:482:LEU:HD12	1.82	0.61
1:A:453:ALA:O	1:A:457:LEU:HG	2.01	0.61
1:D:422:VAL:O	1:D:426:ARG:HG3	2.00	0.61
1:B:330:LEU:HD23	1:B:377:MET:CE	2.31	0.60
1:B:233:ARG:HH21	1:B:439:ARG:HD3	1.65	0.60
1:B:300:ARG:NH2	2:B:633:HOH:O	2.27	0.60
1:C:314:HIS:CD2	1:C:421:ARG:HH11	2.16	0.60
1:D:129:GLY:O	1:D:132:LYS:HD2	2.01	0.59
1:A:351:THR:HB	1:A:352:PRO:HD2	1.83	0.59
1:B:3:ARG:O	1:B:49:HIS:HE1	1.85	0.59
1:D:326:ARG:HD2	1:D:380:GLU:OE2	2.02	0.59
1:B:150:CYS:O	2:B:639:HOH:O	2.15	0.59
1:B:326:ARG:HD2	1:B:380:GLU:OE2	2.03	0.59
1:C:290:HIS:HD2	2:C:655:HOH:O	1.86	0.59
1:C:458:THR:HG22	1:C:459:MET:CE	2.32	0.58
1:C:470:TYR:HB3	2:C:601:HOH:O	2.03	0.58
1:C:179:ASN:ND2	1:C:181:LYS:H	1.98	0.58
1:D:124:THR:HB	1:D:272:LEU:H	1.67	0.58
1:B:31:HIS:HB3	1:B:34:THR:OG1	2.04	0.58
1:D:440:LEU:O	1:D:444:GLU:HG2	2.04	0.58
1:B:351:THR:HB	1:B:352:PRO:CD	2.33	0.58
1:D:67:ARG:HH22	1:D:145:LYS:HE3	1.68	0.58
1:D:24:LYS:NZ	1:D:24:LYS:HB3	2.19	0.58
1:B:134:LYS:CD	1:B:134:LYS:H	2.17	0.57
1:A:390:LEU:H	1:A:390:LEU:CD2	2.16	0.57
1:C:360:ILE:HA	1:C:363:TYR:CE2	2.38	0.57
1:D:452:GLN:HG2	1:D:453:ALA:N	2.18	0.57
1:C:2:ILE:HG13	1:C:147:ASP:OD2	2.05	0.57
1:D:177:GLU:HG2	1:D:178:LYS:N	2.19	0.57
1:B:31:HIS:CD2	1:B:33:LYS:H	2.23	0.57
1:A:273:GLY:O	1:A:275:PRO:HD3	2.03	0.57
1:D:298:SER:HB3	1:D:337:MET:HB3	1.85	0.57
1:A:37:TYR:CZ	1:A:252:GLU:HG3	2.40	0.57
1:C:390:LEU:H	1:C:390:LEU:CD2	2.17	0.56
1:C:472:THR:HA	1:C:475:ARG:NH2	2.20	0.56
1:C:78:GLU:HG2	1:C:81:ARG:NH2	2.20	0.56
1:A:390:LEU:H	1:A:390:LEU:HD23	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:HIS:HD2	1:B:33:LYS:HB2	1.70	0.56
1:A:76:ASP:OD1	1:A:79:GLN:HG3	2.06	0.56
1:D:294:LEU:HD13	1:D:341:ILE:HG13	1.88	0.56
1:B:211:ILE:HD12	1:B:212:PRO:CD	2.35	0.56
1:B:31:HIS:CD2	1:B:33:LYS:HB2	2.40	0.56
1:D:278:ALA:O	1:D:282:PRO:HG2	2.06	0.56
1:D:393:ALA:HB2	1:D:407:PHE:CE2	2.41	0.55
1:C:472:THR:HA	1:C:475:ARG:CZ	2.36	0.55
1:D:142:PHE:CE1	1:D:146:HIS:CD2	2.94	0.55
1:B:394:SER:OG	1:B:397:GLN:HG3	2.06	0.55
1:D:337:MET:HE2	1:D:369:LEU:HD12	1.86	0.55
1:A:314:HIS:CD2	1:A:421:ARG:HH11	2.21	0.55
1:C:302:GLU:HB2	1:C:334:HIS:CE1	2.42	0.55
1:D:176:ILE:HD11	1:D:184:VAL:HG11	1.87	0.55
1:C:224:ASN:C	1:C:224:ASN:HD22	2.10	0.55
1:D:413:GLY:HA3	2:D:602:HOH:O	2.07	0.55
1:C:481:GLU:CG	1:C:482:LEU:H	2.15	0.55
1:D:158:GLN:O	1:D:158:GLN:HG3	2.07	0.55
1:A:105:VAL:HG22	1:A:292:HIS:CE1	2.42	0.55
1:C:458:THR:HG22	1:C:459:MET:HE2	1.89	0.54
1:C:239:GLU:O	1:C:241:PRO:HD3	2.07	0.54
1:B:134:LYS:H	1:B:134:LYS:HD2	1.71	0.54
1:B:31:HIS:HD2	1:B:33:LYS:H	1.54	0.54
1:C:76:ASP:OD1	1:C:79:GLN:HG3	2.07	0.54
1:B:16:ARG:O	2:B:655:HOH:O	2.18	0.54
1:C:88:GLU:HA	1:C:91:MET:CE	2.38	0.54
1:D:84:ARG:HG3	1:D:295:ILE:HD13	1.90	0.54
1:A:31:HIS:CD2	1:A:33:LYS:H	2.23	0.54
1:A:314:HIS:CD2	1:A:421:ARG:HB2	2.43	0.54
1:B:134:LYS:HD2	1:B:134:LYS:N	2.22	0.54
1:D:109:VAL:HG13	1:D:285:VAL:HG12	1.90	0.54
1:B:119:PRO:HA	1:B:122:THR:HB	1.88	0.54
1:C:466:ALA:O	1:C:475:ARG:NH1	2.40	0.54
1:C:452:GLN:O	1:C:456:MET:HG2	2.08	0.53
1:D:31:HIS:HD2	1:D:33:LYS:HB2	1.72	0.53
1:D:208:ARG:HG3	1:D:208:ARG:HH11	1.72	0.53
1:C:351:THR:HB	1:C:352:PRO:HD2	1.91	0.53
1:D:273:GLY:O	1:D:275:PRO:HD3	2.08	0.53
1:D:194:GLY:HA3	2:D:626:HOH:O	2.09	0.53
1:A:439:ARG:HH21	1:A:440:LEU:HD13	1.73	0.53
1:C:88:GLU:HA	1:C:91:MET:HE2	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:50:ARG:HB3	1:D:52:ASP:OD2	2.09	0.53
1:A:481:GLU:HG2	1:A:482:LEU:H	1.73	0.52
1:D:61:ASP:HB2	1:D:67:ARG:HG3	1.90	0.52
1:A:438:ASN:O	1:A:441:PHE:HB3	2.09	0.52
1:B:246:LEU:HD12	1:B:431:LEU:HD13	1.92	0.52
1:C:105:VAL:HG22	1:C:292:HIS:CE1	2.44	0.52
1:A:78:GLU:H	1:A:78:GLU:CD	2.13	0.52
1:C:109:VAL:HG22	1:C:285:VAL:O	2.10	0.52
1:D:439:ARG:HH21	1:D:440:LEU:HD23	1.75	0.52
1:B:274:ASN:HD22	1:B:276:GLU:H	1.56	0.51
1:C:246:LEU:HD12	1:C:431:LEU:HD13	1.92	0.51
1:D:207:PHE:HB3	1:D:279:LYS:HB3	1.92	0.51
1:A:274:ASN:OD1	1:A:276:GLU:HB2	2.11	0.51
1:A:359:ASN:HD21	1:A:361:LEU:HB3	1.76	0.51
1:B:178:LYS:HB3	1:B:183:ILE:HG22	1.92	0.51
1:C:78:GLU:HG2	1:C:81:ARG:HH21	1.75	0.51
1:C:453:ALA:O	1:C:457:LEU:HG	2.09	0.51
1:A:19:TRP:NE1	1:A:24:LYS:HE2	2.25	0.51
1:A:265:PRO:HG2	1:A:268:HIS:HD2	1.75	0.51
1:C:359:ASN:HD22	1:C:362:ILE:H	1.59	0.51
1:D:56:VAL:O	1:D:69:MET:HG3	2.10	0.51
1:D:482:LEU:N	1:D:482:LEU:CD2	2.67	0.51
1:A:292:HIS:HE1	2:A:688:HOH:O	1.92	0.51
1:B:61:ASP:OD2	1:B:67:ARG:HD2	2.11	0.51
1:D:450:PRO:HD2	1:D:453:ALA:CB	2.41	0.50
1:A:61:ASP:HB3	1:A:63:ASP:OD2	2.10	0.50
1:B:237:VAL:CG1	1:B:250:GLY:HA2	2.41	0.50
1:C:224:ASN:ND2	2:C:706:HOH:O	2.36	0.50
1:C:31:HIS:CD2	1:C:32:PRO:HD2	2.47	0.50
1:A:452:GLN:HE22	1:A:456:MET:CE	2.25	0.50
1:B:350:HIS:HD2	1:B:354:GLY:O	1.95	0.50
1:C:159:MET:SD	1:C:208:ARG:NH1	2.85	0.50
1:C:155:VAL:HG21	1:C:191:ILE:HB	1.94	0.50
1:D:276:GLU:O	1:D:279:LYS:HG2	2.11	0.50
1:B:17:ASN:OD1	1:B:228:VAL:O	2.30	0.50
1:B:274:ASN:C	1:B:274:ASN:ND2	2.64	0.50
1:C:31:HIS:CD2	1:C:33:LYS:H	2.25	0.50
1:B:212:PRO:HB2	1:B:214:ASP:OD1	2.12	0.49
1:B:369:LEU:C	1:B:369:LEU:CD2	2.81	0.49
1:D:336:ALA:O	1:D:340:HIS:HD2	1.95	0.49
1:A:390:LEU:HD23	1:A:390:LEU:N	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:208:ARG:HG2	1:C:211:ILE:HB	1.95	0.49
1:D:84:ARG:HG3	1:D:295:ILE:CD1	2.42	0.49
1:A:120:TRP:CZ2	1:A:281:TYR:HB3	2.46	0.49
1:A:27:ASN:C	1:A:27:ASN:HD22	2.15	0.49
1:A:159:MET:CE	1:A:208:ARG:HD3	2.43	0.49
1:C:91:MET:HE1	1:C:295:ILE:HG22	1.93	0.49
1:D:360:ILE:HA	1:D:363:TYR:CE2	2.48	0.49
1:B:131:VAL:HG22	1:B:269:VAL:HG22	1.95	0.49
1:C:314:HIS:HD2	1:C:421:ARG:HD3	1.75	0.49
1:A:76:ASP:OD2	2:A:681:HOH:O	2.20	0.49
1:A:294:LEU:HA	1:A:297:GLN:NE2	2.28	0.49
1:D:41:HIS:O	1:D:44:PHE:HB3	2.13	0.49
1:A:216:ILE:HG22	1:A:216:ILE:O	2.11	0.49
1:C:390:LEU:HG	1:C:391:ILE:CG1	2.35	0.49
1:D:67:ARG:NH2	1:D:145:LYS:HE3	2.28	0.49
1:D:152:PRO:HA	1:D:202:HIS:HB3	1.95	0.49
1:B:84:ARG:NH1	1:B:345:GLU:OE2	2.46	0.48
1:C:139:PHE:O	1:C:142:PHE:HB3	2.13	0.48
1:D:154:PHE:HA	1:D:204:GLY:O	2.12	0.48
1:A:233:ARG:NH2	1:A:439:ARG:HG3	2.28	0.48
1:C:367:ARG:NH2	1:C:449:THR:O	2.45	0.48
1:A:471:GLY:O	1:A:475:ARG:HG3	2.14	0.48
1:A:91:MET:HE2	2:A:688:HOH:O	2.13	0.48
1:D:57:MET:HE3	1:D:86:TYR:HE1	1.78	0.48
1:C:23:GLU:OE2	1:C:31:HIS:HE1	1.97	0.48
1:D:31:HIS:CD2	1:D:33:LYS:HB2	2.49	0.48
1:D:458:THR:HA	1:D:461:ARG:HH11	1.78	0.47
1:B:440:LEU:O	1:B:444:GLU:HG2	2.13	0.47
1:C:224:ASN:C	1:C:224:ASN:ND2	2.67	0.47
1:D:104:ASP:O	2:D:608:HOH:O	2.20	0.47
1:A:117:PRO:HB2	1:A:133:ALA:HB1	1.96	0.47
1:B:122:THR:HG22	1:B:123:GLN:NE2	2.30	0.47
1:C:444:GLU:OE2	1:C:444:GLU:HA	2.14	0.47
1:B:466:ALA:O	1:B:475:ARG:NH1	2.48	0.47
1:D:121:GLU:O	1:D:124:THR:HG23	2.14	0.47
1:D:124:THR:O	1:D:125:ILE:HD13	2.14	0.47
1:B:81:ARG:NH1	1:B:345:GLU:OE2	2.42	0.47
1:A:145:LYS:HD3	1:A:146:HIS:CE1	2.49	0.47
1:B:438:ASN:O	1:B:441:PHE:HB3	2.14	0.47
1:D:130:LYS:HB3	1:D:266:TRP:CD2	2.49	0.47
1:A:318:ASN:H	1:A:318:ASN:ND2	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:ALA:HB3	1:B:98:SER:HB2	1.96	0.47
1:D:439:ARG:HD3	1:D:439:ARG:C	2.35	0.47
1:A:156:ASP:OD2	1:A:205:VAL:HG23	2.15	0.47
1:D:290:HIS:O	1:D:367:ARG:HD3	2.15	0.47
1:B:2:ILE:HG22	1:B:49:HIS:CE1	2.49	0.46
1:C:124:THR:HG21	1:C:131:VAL:HB	1.97	0.46
1:C:136:ILE:HG12	1:C:270:PHE:CD1	2.50	0.46
1:D:119:PRO:O	1:D:123:GLN:CG	2.63	0.46
1:D:146:HIS:HB2	1:D:148:LEU:CD1	2.44	0.46
1:A:27:ASN:ND2	1:A:29:ALA:H	2.13	0.46
1:B:127:ALA:O	1:B:128:GLU:HB2	2.15	0.46
1:C:74:HIS:CD2	1:C:356:TYR:O	2.68	0.46
1:B:290:HIS:O	1:B:367:ARG:HD3	2.16	0.46
1:C:179:ASN:ND2	1:C:182:GLY:H	2.14	0.46
1:C:333:PHE:CD2	1:C:377:MET:HG3	2.51	0.46
1:D:26:ASP:CG	1:D:27:ASN:H	2.15	0.46
1:D:91:MET:CE	1:D:292:HIS:CE1	2.97	0.46
1:D:244:HIS:CD2	1:D:430:ASP:OD2	2.63	0.46
1:C:114:ILE:HG22	1:C:115:ASP:N	2.30	0.46
1:D:420:ASP:O	1:D:424:ILE:HG13	2.16	0.46
1:D:31:HIS:CD2	1:D:33:LYS:H	2.34	0.46
1:D:119:PRO:O	1:D:123:GLN:HG2	2.16	0.46
1:D:271:HIS:CE1	1:D:274:ASN:HB3	2.50	0.46
1:B:135:ASN:HB3	1:B:267:SER:O	2.16	0.46
1:B:351:THR:HB	1:B:352:PRO:HD2	1.98	0.46
1:D:393:ALA:HB2	1:D:407:PHE:CZ	2.50	0.46
1:A:114:ILE:HD11	1:A:140:VAL:HG11	1.97	0.46
1:B:201:ILE:O	1:B:220:ALA:HA	2.16	0.46
1:C:319:LYS:HA	1:C:319:LYS:HD3	1.65	0.46
1:A:31:HIS:CD2	1:A:32:PRO:HD2	2.51	0.46
1:B:130:LYS:HD2	1:B:266:TRP:CD1	2.51	0.46
1:D:266:TRP:HA	1:D:269:VAL:CG1	2.44	0.46
1:D:450:PRO:HD2	1:D:453:ALA:HB3	1.97	0.46
1:A:337:MET:CE	1:A:370:TYR:HB2	2.41	0.46
1:C:227:GLY:HA3	1:C:262:VAL:HG22	1.97	0.46
1:D:338:LEU:C	1:D:338:LEU:HD23	2.36	0.46
1:A:290:HIS:HD2	2:A:615:HOH:O	1.98	0.45
1:B:100:PRO:HD2	1:B:252:GLU:OE2	2.16	0.45
1:C:219:ALA:HB1	1:C:268:HIS:O	2.16	0.45
1:D:6:LYS:O	1:D:10:GLU:HG3	2.16	0.45
1:D:438:ASN:O	1:D:441:PHE:HB3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:314:HIS:HE1	1:A:417:GLN:O	2.00	0.45
1:B:74:HIS:HE1	1:B:108:TYR:OH	1.99	0.45
1:B:78:GLU:H	1:B:78:GLU:HG3	1.62	0.45
1:B:98:SER:O	1:B:440:LEU:HD21	2.17	0.45
1:B:122:THR:HG22	1:B:123:GLN:HE21	1.82	0.45
1:B:422:VAL:O	1:B:426:ARG:HG3	2.16	0.45
1:C:155:VAL:CG2	1:C:191:ILE:HB	2.46	0.45
1:D:31:HIS:HD2	1:D:33:LYS:H	1.65	0.45
1:A:149:ASN:ND2	1:A:149:ASN:C	2.69	0.45
1:D:71:TRP:CE2	1:D:144:LYS:HA	2.51	0.45
1:D:83:LYS:O	1:D:86:TYR:HB3	2.17	0.45
1:D:398:TRP:CZ3	1:D:407:PHE:HB3	2.52	0.45
1:D:75:TYR:CD1	1:D:355:ALA:HA	2.51	0.45
1:D:376:GLN:NE2	2:D:606:HOH:O	2.50	0.45
1:A:436:TRP:O	1:A:440:LEU:HD22	2.16	0.45
1:C:109:VAL:HG13	1:C:285:VAL:HG13	1.97	0.45
1:A:37:TYR:OH	1:A:252:GLU:HG3	2.17	0.45
1:B:118:SER:N	1:B:119:PRO:HD2	2.31	0.45
1:B:75:TYR:H	1:B:79:GLN:HE21	1.65	0.45
1:B:134:LYS:CD	1:B:134:LYS:N	2.78	0.45
1:D:137:VAL:HG12	1:D:141:ASN:ND2	2.32	0.45
1:A:482:LEU:O	1:A:483:THR:HB	2.16	0.45
1:B:341:ILE:O	1:B:345:GLU:HG3	2.17	0.45
1:C:319:LYS:HB2	2:C:636:HOH:O	2.17	0.45
1:B:149:ASN:ND2	1:B:198:ALA:HA	2.33	0.44
1:B:142:PHE:CE2	1:B:146:HIS:CD2	3.05	0.44
1:B:393:ALA:HB2	1:B:407:PHE:CE2	2.52	0.44
1:B:314:HIS:HE1	1:B:417:GLN:O	2.00	0.44
1:C:337:MET:HE2	1:C:366:GLY:O	2.17	0.44
1:A:176:ILE:HD11	1:A:184:VAL:HG12	1.99	0.44
1:B:186:SER:HA	1:B:259:PHE:O	2.17	0.44
1:C:154:PHE:HA	1:C:204:GLY:O	2.17	0.44
1:B:74:HIS:HA	1:B:79:GLN:HE21	1.82	0.44
1:B:201:ILE:HG22	1:B:203:ILE:HD13	2.00	0.44
1:A:351:THR:HB	1:A:352:PRO:CD	2.47	0.44
1:B:264:ILE:HA	1:B:265:PRO:HD2	1.86	0.44
1:B:326:ARG:O	1:B:329:LYS:HB3	2.18	0.44
1:C:216:ILE:CD1	1:C:275:PRO:HG3	2.48	0.44
1:A:139:PHE:O	1:A:142:PHE:HB3	2.18	0.44
1:B:233:ARG:NH2	1:B:439:ARG:HD3	2.32	0.44
1:D:203:ILE:O	1:D:218:PHE:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:208:ARG:HG3	1:D:208:ARG:NH1	2.32	0.44
1:D:327:VAL:O	1:D:331:ILE:HG12	2.18	0.44
1:D:91:MET:HE1	1:D:296:ARG:HB2	1.99	0.44
1:D:281:TYR:N	1:D:282:PRO:CD	2.80	0.44
1:A:201:ILE:HG22	1:A:203:ILE:HD13	1.98	0.43
1:A:225:THR:HG21	1:A:262:VAL:HG13	2.00	0.43
1:A:452:GLN:HE21	1:A:452:GLN:HB2	1.58	0.43
1:B:413:GLY:HA3	2:B:606:HOH:O	2.17	0.43
1:C:297:GLN:HE22	1:C:367:ARG:HA	1.82	0.43
1:D:24:LYS:HB3	1:D:24:LYS:HZ2	1.83	0.43
1:A:80:LEU:CD1	1:A:345:GLU:HG2	2.47	0.43
1:B:77:LYS:O	1:B:81:ARG:HB2	2.18	0.43
1:B:98:SER:O	1:B:100:PRO:HD3	2.18	0.43
1:D:62:LYS:C	1:D:64:GLY:H	2.22	0.43
1:C:390:LEU:HD23	1:C:390:LEU:N	2.31	0.43
1:D:40:ARG:HD2	1:D:249:GLN:HB3	2.00	0.43
1:D:91:MET:HE2	1:D:296:ARG:HG3	1.99	0.43
1:A:297:GLN:HG2	1:A:370:TYR:CD1	2.54	0.43
1:D:120:TRP:CH2	1:D:281:TYR:HB3	2.54	0.43
1:D:440:LEU:H	1:D:440:LEU:HG	1.73	0.43
1:A:176:ILE:HD11	1:A:184:VAL:CG1	2.49	0.43
1:B:69:MET:HB2	1:B:86:TYR:CD1	2.54	0.43
1:B:122:THR:CG2	1:B:123:GLN:HE21	2.31	0.43
1:C:71:TRP:CE2	1:C:144:LYS:HA	2.54	0.43
1:C:74:HIS:HD2	1:C:356:TYR:O	2.01	0.43
1:C:130:LYS:HB3	1:C:266:TRP:CD2	2.54	0.43
1:C:33:LYS:HD2	1:C:33:LYS:HA	1.82	0.43
1:A:71:TRP:CE2	1:A:144:LYS:HA	2.53	0.43
1:D:120:TRP:CZ2	1:D:281:TYR:HB3	2.54	0.43
1:D:53:LEU:HB3	1:D:57:MET:HG2	2.00	0.43
1:A:54:GLN:HE21	1:A:66:ARG:NH2	2.16	0.42
1:C:86:TYR:CE1	1:C:90:ILE:HD11	2.54	0.42
1:A:27:ASN:HD22	1:A:28:VAL:N	2.17	0.42
1:A:124:THR:HG21	1:A:131:VAL:HB	1.99	0.42
1:A:224:ASN:C	1:A:224:ASN:ND2	2.72	0.42
1:B:66:ARG:C	1:B:67:ARG:HG2	2.39	0.42
1:B:127:ALA:O	1:B:130:LYS:HB2	2.19	0.42
1:B:235:SER:OG	1:B:237:VAL:HG22	2.19	0.42
1:A:62:LYS:HG2	1:A:63:ASP:H	1.80	0.42
1:B:61:ASP:OD1	1:B:61:ASP:C	2.58	0.42
1:B:84:ARG:HD3	1:B:345:GLU:OE1	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:66:ARG:HG3	1:D:66:ARG:HH11	1.84	0.42
1:D:225:THR:O	1:D:228:VAL:HB	2.19	0.42
1:D:271:HIS:NE2	1:D:274:ASN:HB3	2.35	0.42
1:A:276:GLU:OE1	1:A:276:GLU:HA	2.20	0.42
1:B:294:LEU:CD2	1:B:337:MET:HG3	2.49	0.42
1:C:271:HIS:HE1	1:C:274:ASN:H	1.67	0.42
1:D:309:ILE:O	1:D:313:GLU:HG3	2.20	0.42
1:A:360:ILE:HA	1:A:363:TYR:CE2	2.54	0.42
1:D:2:ILE:HD11	1:D:145:LYS:HE3	2.02	0.42
1:B:233:ARG:HG2	1:B:254:ASP:O	2.19	0.42
1:B:319:LYS:HD3	1:B:319:LYS:HA	1.82	0.42
1:C:246:LEU:CD1	1:C:431:LEU:HD13	2.49	0.42
1:C:56:VAL:O	1:C:69:MET:HG3	2.19	0.42
1:C:337:MET:CE	1:C:366:GLY:O	2.68	0.42
1:D:173:LEU:O	1:D:215:GLN:HG2	2.20	0.42
1:A:466:ALA:O	1:A:475:ARG:NH1	2.53	0.41
1:C:438:ASN:O	1:C:441:PHE:HB3	2.20	0.41
1:D:37:TYR:CZ	1:D:252:GLU:HG3	2.56	0.41
1:D:281:TYR:O	1:D:285:VAL:HG23	2.20	0.41
1:B:294:LEU:HD23	1:B:294:LEU:HA	1.86	0.41
1:A:54:GLN:NE2	1:A:66:ARG:NH2	2.68	0.41
1:B:6:LYS:HE2	1:B:10:GLU:OE2	2.20	0.41
1:B:120:TRP:O	1:B:271:HIS:HD2	2.02	0.41
1:B:152:PRO:HA	1:B:202:HIS:HB3	2.01	0.41
1:C:289:LEU:HD13	1:C:289:LEU:O	2.20	0.41
1:C:463:GLU:HA	1:C:469:PRO:HD2	2.02	0.41
1:D:129:GLY:O	1:D:132:LYS:CD	2.68	0.41
1:D:37:TYR:OH	1:D:252:GLU:HG3	2.21	0.41
1:D:120:TRP:O	1:D:124:THR:HG22	2.20	0.41
1:B:83:LYS:O	1:B:86:TYR:HB3	2.20	0.41
1:D:221:THR:HG22	1:D:222:PRO:HD2	2.02	0.41
1:A:154:PHE:CE2	1:A:204:GLY:HA3	2.56	0.41
1:B:405:GLY:O	1:B:409:ARG:HG3	2.21	0.41
1:C:31:HIS:CD2	1:C:33:LYS:HB2	2.56	0.41
1:D:118:SER:N	1:D:119:PRO:CD	2.83	0.41
1:A:54:GLN:NE2	1:A:66:ARG:CZ	2.84	0.41
1:A:320:ILE:HA	1:A:321:PRO:HD3	1.94	0.41
1:B:225:THR:HA	1:B:226:PRO:HD3	1.92	0.41
1:B:322:ALA:O	1:B:326:ARG:HG3	2.21	0.41
1:D:221:THR:HG22	1:D:225:THR:OG1	2.21	0.41
1:D:326:ARG:HA	1:D:329:LYS:HE2	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:338:LEU:HD23	1:D:338:LEU:O	2.21	0.40
1:A:278:ALA:O	1:A:282:PRO:HG2	2.21	0.40
1:A:405:GLY:O	1:A:409:ARG:HG3	2.21	0.40
1:B:384:LEU:HD12	1:B:384:LEU:HA	1.97	0.40
1:B:407:PHE:HA	1:B:410:MET:HE2	2.03	0.40
1:D:33:LYS:HD3	1:D:33:LYS:HA	1.87	0.40
1:D:91:MET:HE2	1:D:296:ARG:CG	2.52	0.40
1:D:207:PHE:CB	1:D:279:LYS:HB3	2.50	0.40
1:B:373:ASN:O	1:B:377:MET:HG2	2.22	0.40
1:C:26:ASP:OD1	1:C:27:ASN:N	2.53	0.40
1:A:302:GLU:HB2	1:A:334:HIS:CE1	2.57	0.40
1:B:70:GLN:HB2	1:B:104:ASP:HA	2.02	0.40
1:B:207:PHE:CG	1:B:208:ARG:N	2.90	0.40
1:C:109:VAL:HG13	1:C:285:VAL:CG1	2.51	0.40
1:D:118:SER:HB3	1:D:119:PRO:HD3	2.03	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	467/517 (90%)	446 (96%)	20 (4%)	1 (0%)	47 68
1	B	467/517 (90%)	443 (95%)	21 (4%)	3 (1%)	25 43
1	C	467/517 (90%)	453 (97%)	14 (3%)	0	100 100
1	D	467/517 (90%)	442 (95%)	21 (4%)	4 (1%)	17 31
All	All	1868/2068 (90%)	1784 (96%)	76 (4%)	8 (0%)	34 54

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	128	GLU
1	A	482	LEU
1	B	172	GLY
1	D	157	PRO
1	B	127	ALA
1	B	390	LEU
1	D	63	ASP
1	D	280	LEU

### 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	399/435 (92%)	366 (92%)	33 (8%)	11	22
1	B	398/435 (92%)	370 (93%)	28 (7%)	15	29
1	C	399/435 (92%)	370 (93%)	29 (7%)	14	27
1	D	399/435 (92%)	367 (92%)	32 (8%)	12	23
All	All	1595/1740 (92%)	1473 (92%)	122 (8%)	13	25

All (122) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	GLU
1	A	12	LEU
1	A	27	ASN
1	A	50	ARG
1	A	62	LYS
1	A	67	ARG
1	A	78	GLU
1	A	80	LEU
1	A	108	TYR
1	A	109	VAL
1	A	149	ASN
1	A	179	ASN
1	A	186	SER
1	A	224	ASN

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	239	GLU
1	A	260	ASP
1	A	284	ARG
1	A	318	ASN
1	A	326	ARG
1	A	330	LEU
1	A	338	LEU
1	A	359	ASN
1	A	390	LEU
1	A	417	GLN
1	A	421	ARG
1	A	431	LEU
1	A	435	ASP
1	A	440	LEU
1	A	452	GLN
1	A	456	MET
1	A	457	LEU
1	A	473	LEU
1	A	482	LEU
1	B	67	ARG
1	B	76	ASP
1	B	81	ARG
1	B	108	TYR
1	B	148	LEU
1	B	174	ARG
1	B	177	GLU
1	B	179	ASN
1	B	208	ARG
1	B	224	ASN
1	B	237	VAL
1	B	269	VAL
1	B	274	ASN
1	B	280	LEU
1	B	289	LEU
1	B	294	LEU
1	B	325	THR
1	B	384	LEU
1	B	408	GLU
1	B	412	ASN
1	B	421	ARG
1	B	431	LEU
1	B	435	ASP

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	B	439	ARG
1	B	457	LEU
1	B	458	THR
1	B	463	GLU
1	B	473	LEU
1	C	12	LEU
1	C	33	LYS
1	C	53	LEU
1	C	56	VAL
1	C	61	ASP
1	C	63	ASP
1	C	66	ARG
1	C	80	LEU
1	C	82	ARG
1	C	108	TYR
1	C	109	VAL
1	C	179	ASN
1	C	202	HIS
1	C	214	ASP
1	C	224	ASN
1	C	260	ASP
1	C	289	LEU
1	C	303	LEU
1	C	329	LYS
1	C	330	LEU
1	C	337	MET
1	C	359	ASN
1	C	361	LEU
1	C	369	LEU
1	C	390	LEU
1	C	421	ARG
1	C	431	LEU
1	C	440	LEU
1	C	463	GLU
1	D	2	ILE
1	D	6	LYS
1	D	62	LYS
1	D	67	ARG
1	D	108	TYR
1	D	109	VAL
1	D	124	THR
1	D	145	LYS

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Mol	Chain	Res	Type
1	D	158	GLN
1	D	174	ARG
1	D	180	ASP
1	D	188	VAL
1	D	191	ILE
1	D	221	THR
1	D	294	LEU
1	D	300	ARG
1	D	325	THR
1	D	359	ASN
1	D	369	LEU
1	D	384	LEU
1	D	388	SER
1	D	391	ILE
1	D	409	ARG
1	D	421	ARG
1	D	431	LEU
1	D	440	LEU
1	D	452	GLN
1	D	457	LEU
1	D	458	THR
1	D	460	GLN
1	D	482	LEU
1	D	483	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (76) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	31	HIS
1	A	39	GLN
1	A	54	GLN
1	A	149	ASN
1	A	179	ASN
1	A	224	ASN
1	A	268	HIS
1	A	290	HIS
1	A	292	HIS
1	A	297	GLN
1	A	314	HIS
1	A	318	ASN
1	A	334	HIS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	340	HIS
1	A	350	HIS
1	A	359	ASN
1	A	397	GLN
1	A	417	GLN
1	A	452	GLN
1	B	31	HIS
1	B	49	HIS
1	B	74	HIS
1	B	79	GLN
1	B	123	GLN
1	B	146	HIS
1	B	179	ASN
1	B	224	ASN
1	B	244	HIS
1	B	249	GLN
1	B	271	HIS
1	B	274	ASN
1	B	283	GLN
1	B	314	HIS
1	B	340	HIS
1	B	350	HIS
1	B	373	ASN
1	B	376	GLN
1	B	397	GLN
1	B	404	ASN
1	B	417	GLN
1	B	438	ASN
1	C	31	HIS
1	C	74	HIS
1	C	135	ASN
1	C	179	ASN
1	C	224	ASN
1	C	268	HIS
1	C	271	HIS
1	C	290	HIS
1	C	292	HIS
1	C	297	GLN
1	C	314	HIS
1	C	318	ASN
1	C	335	GLN
1	C	340	HIS

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Mol	Chain	Res	Type
1	C	359	ASN
1	C	376	GLN
1	C	397	GLN
1	C	438	ASN
1	D	31	HIS
1	D	39	GLN
1	D	74	HIS
1	D	141	ASN
1	D	146	HIS
1	D	158	GLN
1	D	244	HIS
1	D	249	GLN
1	D	314	HIS
1	D	318	ASN
1	D	334	HIS
1	D	340	HIS
1	D	350	HIS
1	D	359	ASN
1	D	376	GLN
1	D	397	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](#)

There are no ligands in this entry.



## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	471/517 (91%)	-0.68	3 (0%) 89 90	30, 44, 67, 98	0
1	B	471/517 (91%)	-0.51	0 100 100	33, 49, 77, 103	0
1	C	471/517 (91%)	-0.74	1 (0%) 95 95	30, 41, 61, 94	0
1	D	471/517 (91%)	-0.39	6 (1%) 77 79	35, 52, 77, 116	0
All	All	1884/2068 (91%)	-0.58	10 (0%) 91 91	30, 46, 73, 116	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	482	LEU	4.9
1	D	483	THR	4.0
1	A	483	THR	3.3
1	D	126	GLY	3.2
1	D	175	VAL	2.8
1	C	482	LEU	2.8
1	D	127	ALA	2.7
1	D	482	LEU	2.3
1	A	157	PRO	2.1
1	D	159	MET	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

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.