



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:07 PM GMT

PDB ID : 3SY2
Title : Crystal structure of the Salmonella E3 ubiquitin ligase SopA in complex with the human E2 UbcH7
Authors : Diao, J.; Lin, D.Y.; Chen, J.
Deposited on : 2011-07-15
Resolution : 3.27 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

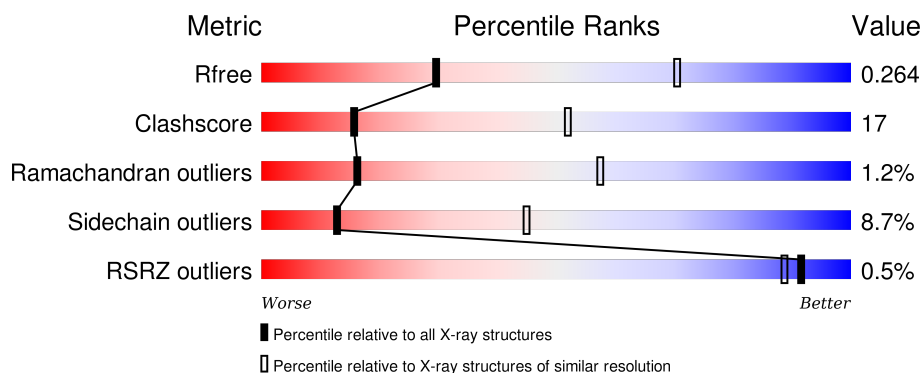
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.27 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1756 (3.36-3.20)
Clashscore	102246	1941 (3.36-3.20)
Ramachandran outliers	100387	1905 (3.36-3.20)
Sidechain outliers	100360	1903 (3.36-3.20)
RSRZ outliers	91569	1764 (3.36-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	621	<div> <div>%</div> <div> <div></div> <div>63%</div> <div>31%</div> <div>• •</div> </div> </div>
1	B	621	<div> <div></div> <div> <div>65%</div> <div>31%</div> <div>• •</div> </div> </div>
2	C	156	<div> <div></div> <div> <div>60%</div> <div>36%</div> <div>• •</div> </div> </div>
2	D	156	<div> <div>%</div> <div> <div></div> <div>58%</div> <div>37%</div> <div>• •</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	1	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11660 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 E3 ubiquitin-protein ligase SopA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	609	Total	C	N	O	S	0	0	0
			4631	2935	794	882	20			
1	B	616	Total	C	N	O	S	0	0	0
			4667	2965	796	886	20			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	162	SER	-	EXPRESSION TAG	UNP Q8ZNR3
A	163	ASN	-	EXPRESSION TAG	UNP Q8ZNR3
A	164	ALA	-	EXPRESSION TAG	UNP Q8ZNR3
B	162	SER	-	EXPRESSION TAG	UNP Q8ZNR3
B	163	ASN	-	EXPRESSION TAG	UNP Q8ZNR3
B	164	ALA	-	EXPRESSION TAG	UNP Q8ZNR3

- Molecule 2 is a protein called Ubiquitin-conjugating enzyme E2 L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	154	Total	C	N	O	S	0	0	0
			1175	753	206	212	4			
2	D	154	Total	C	N	O	S	0	0	0
			1167	751	202	209	5			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	GLY	-	EXPRESSION TAG	UNP P68036
C	0	SER	-	EXPRESSION TAG	UNP P68036
D	-1	GLY	-	EXPRESSION TAG	UNP P68036
D	0	SER	-	EXPRESSION TAG	UNP P68036

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

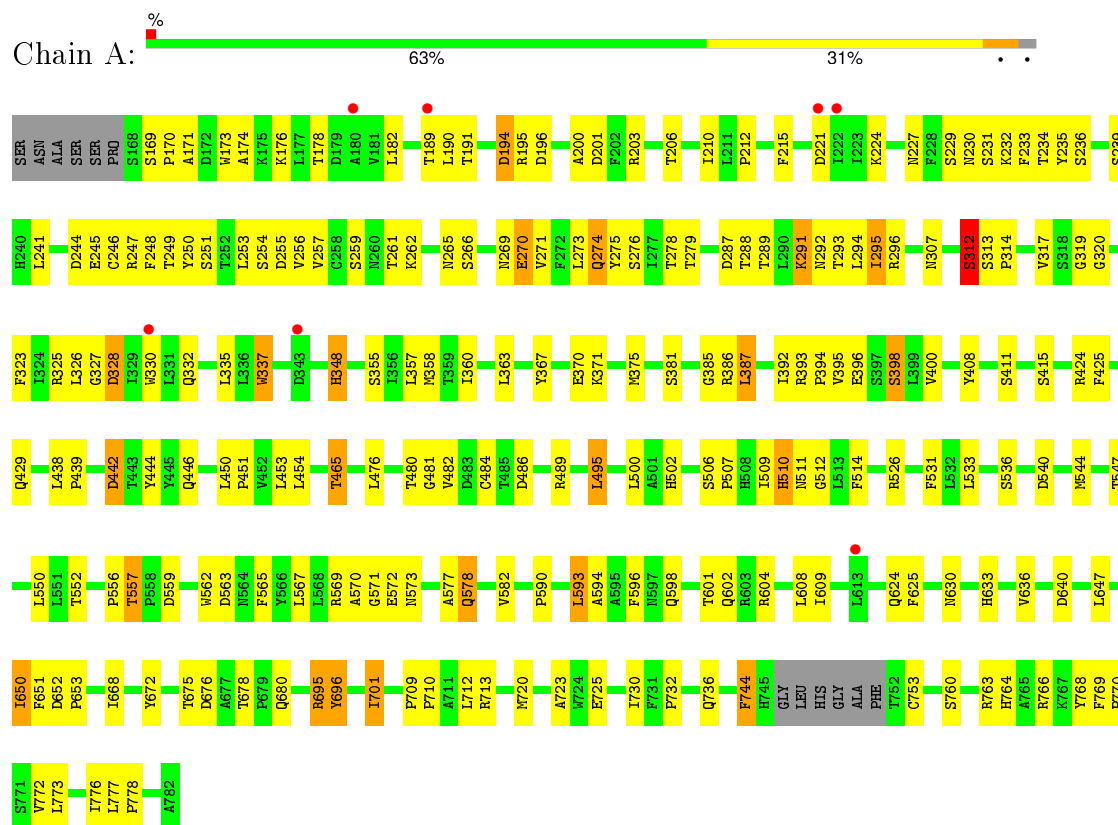
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	8	Total	O	0	0
			8	8		
4	B	6	Total	O	0	0
			6	6		
4	D	1	Total	O	0	0
			1	1		

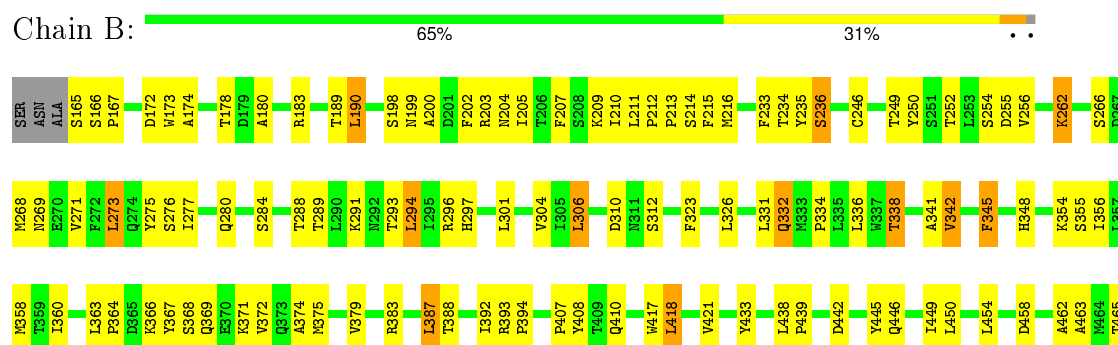
3 Residue-property plots

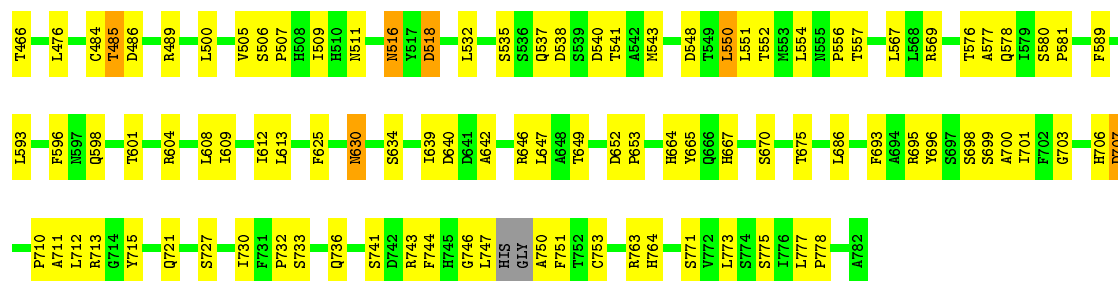
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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: E3 ubiquitin-protein ligase SopA



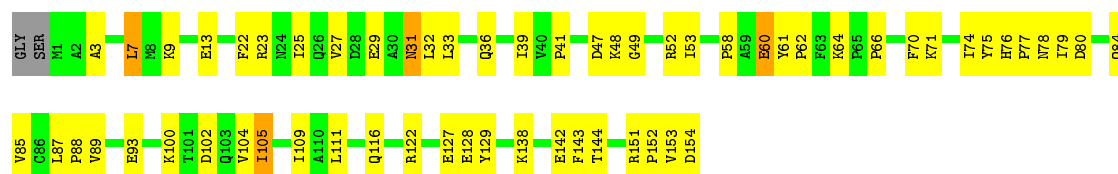
• Molecule 1: E3 ubiquitin-protein ligase SopA





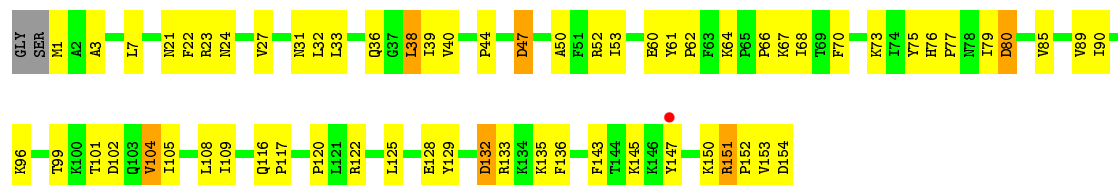
• Molecule 2: Ubiquitin-conjugating enzyme E2 L3

Chain C: 60% 36% 4%



• Molecule 2: Ubiquitin-conjugating enzyme E2 L3

Chain D: 58% 37% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.12Å 118.36Å 241.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.42 – 3.27 45.42 – 3.27	Depositor EDS
% Data completeness (in resolution range)	93.8 (45.42-3.27) 93.7 (45.42-3.27)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 3.25Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.211 , 0.275 0.200 , 0.264	Depositor DCC
R_{free} test set	1570 reflections (4.97%)	DCC
Wilson B-factor (Å ²)	85.2	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 61.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 31621 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11660	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [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	0.28	0/4752	0.46	0/6493
1	B	0.29	0/4790	0.46	0/6548
2	C	0.29	0/1206	0.49	0/1646
2	D	0.30	0/1197	0.54	0/1635
All	All	0.29	0/11945	0.47	0/16322

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	4631	0	4322	160	0
1	B	4667	0	4363	134	0
2	C	1175	0	1107	44	0
2	D	1167	0	1110	50	0
3	A	5	0	0	0	0
4	A	8	0	0	0	0
4	B	6	0	0	0	0
4	D	1	0	0	0	0
All	All	11660	0	10902	385	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 17.

All (385) 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)
2:D:151:ARG:HD3	2:D:152:PRO:HD2	1.45	0.97
2:C:52:ARG:HH12	2:C:71:LYS:HD3	1.37	0.88
1:A:375:MET:HG3	1:A:408:TYR:HE2	1.37	0.88
1:A:274:GLN:HG3	1:A:294:LEU:HD22	1.53	0.87
1:A:593:LEU:HD23	1:A:594:ALA:N	1.89	0.87
2:D:96:LYS:O	2:D:99:THR:HG23	1.76	0.85
2:D:23:ARG:O	2:D:38:LEU:HB3	1.75	0.85
2:C:78:ASN:HD21	2:C:116:GLN:H	1.25	0.84
1:B:438:LEU:HD12	1:B:439:PRO:HD2	1.59	0.84
1:B:174:ALA:HB1	1:B:210:ILE:HD12	1.60	0.84
1:A:375:MET:HG3	1:A:408:TYR:CE2	2.12	0.84
2:C:76:HIS:HD2	2:C:78:ASN:H	1.26	0.83
1:B:387:LEU:H	1:B:387:LEU:HD12	1.42	0.83
1:A:590:PRO:O	1:A:593:LEU:HD22	1.78	0.83
1:A:680:GLN:HG3	1:A:730:ILE:HD11	1.61	0.83
1:B:732:PRO:HD3	1:B:764:HIS:CE1	2.14	0.82
1:B:695:ARG:HH11	1:B:695:ARG:HG3	1.46	0.80
2:D:101:THR:HA	2:D:104:VAL:HG13	1.63	0.79
1:A:244:ASP:OD2	1:A:245:GLU:HG3	1.83	0.78
1:A:593:LEU:HD23	1:A:594:ALA:H	1.44	0.78
2:C:78:ASN:ND2	2:C:111:LEU:HD11	1.99	0.76
2:C:52:ARG:NH1	2:C:71:LYS:HD3	2.00	0.76
1:A:582:VAL:HG22	1:A:596:PHE:CE2	2.20	0.75
2:C:76:HIS:CD2	2:C:78:ASN:H	2.05	0.75
2:D:7:LEU:HD11	2:D:33:LEU:HD23	1.71	0.72
1:B:743:ARG:HD2	1:B:753:CYS:SG	2.29	0.72
1:A:393:ARG:CZ	1:A:763:ARG:HD3	2.20	0.72
1:A:325:ARG:HH21	1:A:330:TRP:HE1	1.38	0.71
1:A:557:THR:HG23	1:A:559:ASP:H	1.56	0.71
1:A:319:GLY:N	1:A:320:GLY:HA3	2.06	0.70
1:B:727:SER:O	1:B:730:ILE:HG12	1.91	0.70
1:A:307:ASN:O	1:A:328:ASP:HB3	1.92	0.69
1:B:233:PHE:O	1:B:236:SER:HB3	1.91	0.69
1:A:438:LEU:HD12	1:A:439:PRO:HD2	1.73	0.69
1:A:190:LEU:HB3	1:A:194:ASP:HB2	1.75	0.68
1:B:277:ILE:HD12	1:B:706:HIS:CD2	2.28	0.68
1:A:540:ASP:HB2	1:A:569:ARG:NH1	2.09	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:3:ALA:O	2:C:7:LEU:HB2	1.94	0.68
1:A:732:PRO:HG2	1:A:736:GLN:HG2	1.76	0.67
1:B:288:THR:HG22	1:B:289:THR:H	1.58	0.67
1:A:500:LEU:HD22	1:A:509:ILE:HD11	1.76	0.67
1:A:190:LEU:HB3	1:A:194:ASP:CB	2.25	0.67
1:A:577:ALA:O	1:A:578:GLN:HB2	1.94	0.67
2:D:22:PHE:HD1	2:D:39:ILE:HG22	1.58	0.67
2:C:79:ILE:HG12	2:C:85:VAL:HG22	1.77	0.66
1:A:573:ASN:ND2	2:D:96:LYS:HB2	2.10	0.66
1:A:251:SER:H	1:A:271:VAL:HG12	1.60	0.66
1:A:732:PRO:HD3	1:A:764:HIS:CE1	2.31	0.65
1:A:292:ASN:HD21	1:A:312:SER:CB	2.09	0.65
1:A:269:ASN:O	1:A:271:VAL:HG13	1.96	0.65
1:A:271:VAL:O	1:A:293:THR:HG23	1.96	0.65
1:B:598:GLN:HA	1:B:601:THR:HB	1.79	0.65
2:D:22:PHE:CD1	2:D:39:ILE:HG22	2.31	0.64
2:C:80:ASP:HB3	2:C:84:GLN:H	1.62	0.64
1:B:165:SER:C	1:B:167:PRO:HD3	2.18	0.64
1:B:306:LEU:HB2	1:B:367:TYR:CE2	2.31	0.64
1:B:198:SER:O	1:B:199:ASN:HB2	1.97	0.64
1:A:174:ALA:HB1	1:A:210:ILE:CD1	2.28	0.64
2:C:61:TYR:CD1	2:C:62:PRO:HA	2.33	0.64
2:D:79:ILE:HG12	2:D:85:VAL:HG22	1.79	0.64
1:A:385:GLY:O	1:A:386:ARG:HB2	1.97	0.63
2:D:117:PRO:O	2:D:120:PRO:HD3	1.99	0.63
1:B:364:PRO:HG2	1:B:367:TYR:CD1	2.34	0.62
2:C:22:PHE:HD1	2:C:39:ILE:HG22	1.63	0.62
1:A:254:SER:O	1:A:255:ASP:HB2	1.99	0.62
1:B:713:ARG:HH12	1:B:746:GLY:HA3	1.63	0.62
2:D:21:ASN:HB3	2:D:109:ILE:HD13	1.81	0.62
1:B:341:ALA:O	1:B:345:PHE:HB2	1.99	0.62
1:B:269:ASN:ND2	1:B:291:LYS:HB3	2.15	0.62
1:A:540:ASP:HB2	1:A:569:ARG:HH12	1.65	0.62
1:B:345:PHE:O	1:B:356:ILE:HG22	2.00	0.62
1:B:166:SER:N	1:B:167:PRO:HD3	2.15	0.61
1:A:695:ARG:HH11	1:A:695:ARG:HG3	1.65	0.61
1:B:695:ARG:HG3	1:B:695:ARG:NH1	2.14	0.61
1:A:348:HIS:HB3	1:A:355:SER:HB3	1.83	0.61
1:B:500:LEU:HA	1:B:505:VAL:HG11	1.82	0.60
2:D:27:VAL:CG2	2:D:32:LEU:HA	2.31	0.60
1:B:476:LEU:HD21	1:B:551:LEU:HD23	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:773:LEU:HD11	1:A:777:LEU:HD12	1.83	0.60
2:D:75:TYR:HB2	2:D:143:PHE:CD1	2.36	0.60
1:B:348:HIS:HB3	1:B:355:SER:HB3	1.83	0.60
1:A:768:TYR:OH	2:C:127:GLU:HG2	2.01	0.60
1:A:246:CYS:H	1:A:266:SER:CB	2.15	0.59
2:C:74:ILE:HG12	2:C:75:TYR:N	2.17	0.59
2:C:58:PRO:HD3	2:C:66:PRO:HA	1.82	0.59
1:A:313:SER:HB2	1:A:314:PRO:HD2	1.84	0.59
2:D:61:TYR:CD1	2:D:62:PRO:HA	2.37	0.59
1:A:169:SER:HB3	1:A:170:PRO:HD2	1.85	0.59
1:B:485:THR:O	1:B:489:ARG:HG3	2.03	0.59
1:A:442:ASP:O	1:A:446:GLN:HG3	2.01	0.59
1:A:234:THR:O	1:A:235:TYR:HB2	2.02	0.59
1:B:334:PRO:HD3	1:B:345:PHE:CZ	2.38	0.58
1:B:634:SER:HB3	1:B:701:ILE:HA	1.85	0.58
1:B:554:LEU:C	1:B:556:PRO:HD3	2.24	0.58
2:D:60:GLU:HG3	2:D:64:LYS:HD2	1.86	0.57
1:A:604:ARG:HG2	1:A:672:TYR:CE2	2.39	0.57
1:A:481:GLY:HA3	1:A:484:CYS:SG	2.45	0.57
1:A:348:HIS:HD1	1:A:358:MET:HG2	1.68	0.57
2:C:78:ASN:ND2	2:C:116:GLN:H	1.98	0.57
1:A:326:LEU:HD11	1:A:363:LEU:HD11	1.87	0.57
1:B:277:ILE:HD12	1:B:706:HIS:CG	2.40	0.56
2:D:31:ASN:O	2:D:32:LEU:HB3	2.05	0.56
1:A:203:ARG:HG3	1:A:232:LYS:O	2.04	0.56
1:A:327:GLY:HA2	1:A:370:GLU:HG2	1.87	0.56
2:C:128:GLU:HG2	2:C:128:GLU:O	2.05	0.56
1:B:256:VAL:O	1:B:275:TYR:HB2	2.05	0.56
1:B:207:PHE:HA	1:B:210:ILE:HG12	1.88	0.56
2:C:74:ILE:HG12	2:C:75:TYR:H	1.70	0.56
2:C:53:ILE:HD11	2:C:70:PHE:CE1	2.41	0.55
2:D:27:VAL:HG21	2:D:32:LEU:HD12	1.89	0.55
1:B:577:ALA:HB1	2:C:93:GLU:HA	1.89	0.55
1:B:275:TYR:CE1	1:B:296:ARG:HD3	2.41	0.55
1:A:625:PHE:HD2	1:A:696:TYR:CZ	2.24	0.55
1:B:732:PRO:HG2	1:B:736:GLN:HG2	1.87	0.55
2:C:9:LYS:HE3	2:C:13:GLU:OE2	2.07	0.55
2:D:61:TYR:CG	2:D:62:PRO:HA	2.42	0.55
1:B:255:ASP:O	1:B:276:SER:HB2	2.07	0.55
1:B:516:ASN:C	1:B:516:ASN:HD22	2.11	0.55
1:A:224:LYS:HA	1:A:244:ASP:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:THR:HG23	1:A:269:ASN:HB2	1.89	0.54
1:A:348:HIS:HB3	1:A:355:SER:CB	2.38	0.54
1:B:446:GLN:HA	1:B:450:LEU:HD12	1.89	0.54
1:B:407:PRO:O	1:B:410:GLN:HG2	2.07	0.54
1:B:393:ARG:N	1:B:394:PRO:HD2	2.22	0.54
2:C:76:HIS:CD2	2:C:77:PRO:HD2	2.43	0.53
1:A:396:GLU:O	1:A:400:VAL:HG22	2.08	0.53
1:B:294:LEU:HD21	1:B:332:GLN:HG3	1.91	0.53
1:A:337:TRP:HZ2	1:A:381:SER:HG	1.56	0.53
1:B:273:LEU:O	1:B:273:LEU:HD23	2.09	0.53
1:A:393:ARG:HA	1:A:396:GLU:HG3	1.91	0.53
1:A:276:SER:C	1:A:278:THR:H	2.12	0.53
2:D:3:ALA:O	2:D:7:LEU:HG	2.08	0.52
1:A:557:THR:CG2	1:A:559:ASP:H	2.21	0.52
1:B:293:THR:HG22	1:B:294:LEU:N	2.24	0.52
2:C:61:TYR:CG	2:C:62:PRO:HA	2.43	0.52
1:B:476:LEU:HD22	1:B:550:LEU:HD13	1.91	0.52
2:D:129:TYR:O	2:D:133:ARG:HG2	2.10	0.52
1:B:535:SER:HB3	1:B:538:ASP:O	2.10	0.52
1:B:609:ILE:HD13	1:B:625:PHE:HB3	1.92	0.52
1:A:348:HIS:ND1	1:A:358:MET:HG2	2.25	0.52
1:B:463:ALA:HA	1:B:466:THR:OG1	2.09	0.52
1:A:476:LEU:HD22	1:A:550:LEU:HD23	1.91	0.52
1:B:207:PHE:HB3	1:B:211:LEU:HD12	1.91	0.52
1:B:167:PRO:HA	1:B:173:TRP:NE1	2.25	0.51
1:B:271:VAL:O	1:B:293:THR:HG23	2.10	0.51
1:B:368:SER:HB2	1:B:407:PRO:HG3	1.92	0.51
1:A:604:ARG:HG2	1:A:672:TYR:CD2	2.46	0.51
1:A:465:THR:HG21	1:A:502:HIS:ND1	2.26	0.51
2:D:75:TYR:HB2	2:D:143:PHE:HD1	1.76	0.51
1:A:392:ILE:HG21	1:A:424:ARG:HH21	1.75	0.51
1:B:249:THR:O	1:B:250:TYR:HB2	2.10	0.51
2:C:22:PHE:CD1	2:C:39:ILE:HG22	2.45	0.51
1:A:730:ILE:O	1:A:764:HIS:HE1	1.94	0.50
2:C:75:TYR:HB2	2:C:143:PHE:CD1	2.46	0.50
2:D:128:GLU:HG2	2:D:136:PHE:HB2	1.92	0.50
1:A:294:LEU:HD23	1:A:295:ILE:N	2.27	0.50
1:B:288:THR:HG22	1:B:289:THR:N	2.25	0.50
1:B:462:ALA:O	1:B:466:THR:HG23	2.11	0.50
1:A:288:THR:HG22	1:A:289:THR:N	2.27	0.50
1:B:203:ARG:O	1:B:204:ASN:HB2	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:PHE:CE2	1:A:221:ASP:HB2	2.47	0.50
1:B:277:ILE:HB	1:B:706:HIS:CE1	2.46	0.50
1:B:516:ASN:ND2	1:B:518:ASP:H	2.09	0.50
1:B:639:ILE:HG21	1:B:707:ASP:HB3	1.93	0.50
1:A:174:ALA:HB1	1:A:210:ILE:HD11	1.93	0.50
1:B:548:ASP:O	1:B:552:THR:HG23	2.11	0.50
2:C:151:ARG:HG3	2:C:152:PRO:HD2	1.94	0.50
1:A:261:THR:HG22	1:A:262:LYS:N	2.26	0.50
2:D:38:LEU:CD2	2:D:50:ALA:HB1	2.42	0.50
1:B:178:THR:HG22	1:B:211:LEU:HD23	1.94	0.50
1:B:375:MET:HG3	1:B:408:TYR:CE2	2.46	0.50
2:D:38:LEU:HD21	2:D:50:ALA:HB1	1.94	0.50
1:A:189:THR:O	1:A:190:LEU:HD23	2.12	0.50
1:B:433:TYR:CG	1:B:438:LEU:HD13	2.47	0.49
2:D:31:ASN:C	2:D:33:LEU:H	2.15	0.49
1:A:174:ALA:O	1:A:178:THR:HG23	2.12	0.49
1:A:732:PRO:CG	1:A:736:GLN:HG2	2.40	0.49
1:A:249:THR:HG22	1:A:250:TYR:CD1	2.47	0.49
1:B:713:ARG:HH12	1:B:746:GLY:CA	2.26	0.49
2:C:41:PRO:O	2:C:151:ARG:NH2	2.42	0.49
1:B:540:ASP:HB2	1:B:569:ARG:NH1	2.27	0.49
2:C:138:LYS:O	2:C:142:GLU:HG3	2.12	0.49
1:A:348:HIS:CD2	1:A:398:SER:HB3	2.48	0.49
1:B:554:LEU:O	1:B:556:PRO:HD3	2.13	0.49
1:A:476:LEU:HA	1:A:531:PHE:HZ	1.78	0.49
2:D:122:ARG:HH11	2:D:125:LEU:HG	1.78	0.49
1:A:249:THR:O	1:A:250:TYR:HB2	2.13	0.49
1:B:234:THR:O	1:B:235:TYR:HB2	2.11	0.49
1:A:766:ARG:O	1:A:770:PRO:HB3	2.12	0.48
2:D:22:PHE:CE2	2:D:105:ILE:HD13	2.48	0.48
1:A:288:THR:CG2	1:A:289:THR:N	2.76	0.48
1:A:652:ASP:N	1:A:653:PRO:CD	2.76	0.48
1:A:582:VAL:HG22	1:A:596:PHE:HE2	1.77	0.48
1:A:276:SER:C	1:A:278:THR:N	2.64	0.48
1:B:387:LEU:H	1:B:387:LEU:CD1	2.21	0.48
2:D:7:LEU:CD1	2:D:33:LEU:HD23	2.42	0.48
1:A:190:LEU:HB2	1:A:195:ARG:HG3	1.94	0.48
1:A:266:SER:O	1:A:288:THR:HG23	2.13	0.48
1:B:664:HIS:HE1	1:B:715:TYR:OH	1.96	0.48
1:A:244:ASP:OD2	1:A:245:GLU:N	2.47	0.48
1:B:178:THR:HG22	1:B:211:LEU:CD2	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:THR:O	1:A:182:LEU:HG	2.13	0.48
2:D:60:GLU:H	2:D:60:GLU:CD	2.16	0.48
1:A:357:LEU:HA	1:A:375:MET:HE1	1.95	0.47
1:A:173:TRP:CZ2	1:A:200:ALA:HA	2.49	0.47
1:B:750:ALA:O	1:B:753:CYS:HB3	2.13	0.47
1:B:699:SER:HA	1:B:703:GLY:O	2.15	0.47
1:B:665:TYR:HE1	1:B:686:LEU:HD11	1.79	0.47
1:A:647:LEU:O	1:A:650:ILE:HG13	2.14	0.47
1:A:772:VAL:O	1:A:776:ILE:HD13	2.15	0.47
2:D:117:PRO:HG3	2:D:129:TYR:CD1	2.50	0.47
1:B:174:ALA:O	1:B:178:THR:HG23	2.14	0.47
2:C:27:VAL:CG1	2:C:32:LEU:HA	2.44	0.47
1:B:418:LEU:HD22	1:B:421:VAL:HB	1.96	0.47
1:A:577:ALA:O	1:A:578:GLN:CB	2.62	0.47
2:D:70:PHE:CE2	2:D:85:VAL:HG21	2.49	0.47
1:A:720:MET:O	1:A:723:ALA:HB3	2.15	0.47
1:B:773:LEU:HD11	1:B:777:LEU:HD12	1.97	0.47
1:B:640:ASP:OD2	1:B:642:ALA:HB3	2.15	0.47
1:B:190:LEU:H	1:B:190:LEU:HD12	1.80	0.47
1:A:253:LEU:HD12	1:A:253:LEU:N	2.30	0.47
1:B:630:ASN:N	1:B:630:ASN:OD1	2.48	0.47
1:A:489:ARG:HG2	1:A:489:ARG:HH11	1.80	0.47
1:A:171:ALA:O	1:A:174:ALA:HB3	2.14	0.47
2:D:23:ARG:NH1	2:D:154:ASP:OD1	2.41	0.47
1:B:348:HIS:HB3	1:B:355:SER:CB	2.45	0.46
1:B:698:SER:C	1:B:700:ALA:H	2.19	0.46
1:B:751:PHE:C	1:B:753:CYS:N	2.69	0.46
1:B:751:PHE:C	1:B:753:CYS:H	2.18	0.46
2:C:74:ILE:HG21	2:C:79:ILE:HD12	1.98	0.46
1:A:273:LEU:HD12	1:A:293:THR:HG21	1.97	0.46
2:C:60:GLU:OE2	2:C:64:LYS:HD3	2.15	0.46
1:A:327:GLY:HA2	1:A:370:GLU:CG	2.44	0.46
1:A:227:ASN:HA	1:A:247:ARG:HB3	1.96	0.46
1:B:746:GLY:HA3	1:B:747:LEU:HA	1.50	0.46
1:A:323:PHE:HD2	1:A:332:GLN:HA	1.81	0.46
1:A:256:VAL:CG1	1:A:257:VAL:N	2.78	0.46
1:A:393:ARG:N	1:A:394:PRO:HD2	2.31	0.46
2:C:87:LEU:CD1	2:C:89:VAL:HG12	2.46	0.46
1:B:332:GLN:H	1:B:332:GLN:CD	2.19	0.46
1:A:510:HIS:O	1:A:512:GLY:N	2.49	0.46
1:B:647:LEU:HB3	1:B:711:ALA:CB	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:38:LEU:HD21	2:D:50:ALA:CB	2.46	0.45
1:A:231:SER:HB2	1:A:233:PHE:CE1	2.51	0.45
1:A:713:ARG:NH1	1:A:744:PHE:O	2.49	0.45
1:A:651:PHE:CZ	1:A:712:LEU:HD23	2.51	0.45
1:A:234:THR:O	1:A:235:TYR:CB	2.64	0.45
2:C:127:GLU:C	2:C:129:TYR:H	2.20	0.45
1:A:563:ASP:HA	1:A:565:PHE:CE1	2.52	0.45
2:D:40:VAL:HG21	2:D:152:PRO:HG3	1.97	0.45
1:A:696:TYR:HD2	1:A:701:ILE:HG21	1.81	0.45
1:B:393:ARG:CZ	1:B:763:ARG:HD2	2.47	0.45
1:A:212:PRO:HG2	1:A:215:PHE:CD1	2.51	0.45
1:B:293:THR:CG2	1:B:294:LEU:N	2.80	0.45
1:B:207:PHE:C	1:B:209:LYS:H	2.20	0.45
1:B:348:HIS:CD2	1:B:358:MET:HE3	2.52	0.45
1:A:604:ARG:HD2	1:A:672:TYR:HA	1.99	0.45
1:A:495:LEU:HA	1:A:495:LEU:HD23	1.77	0.45
1:B:646:ARG:O	1:B:649:THR:HB	2.17	0.45
1:B:268:MET:HG3	1:B:288:THR:HG21	1.99	0.45
1:B:368:SER:O	1:B:372:VAL:HG23	2.17	0.45
1:A:328:ASP:CG	1:A:328:ASP:O	2.54	0.45
2:D:145:LYS:HE2	2:D:145:LYS:HB3	1.84	0.45
1:A:506:SER:N	1:A:507:PRO:CD	2.80	0.44
2:D:24:ASN:O	2:D:24:ASN:OD1	2.35	0.44
2:D:68:ILE:HD13	2:D:108:LEU:HD12	1.99	0.44
1:B:306:LEU:HD22	1:B:367:TYR:CD2	2.52	0.44
1:B:506:SER:N	1:B:507:PRO:CD	2.81	0.44
1:A:236:SER:O	1:A:256:VAL:HG13	2.18	0.44
1:B:543:MET:HE3	1:B:589:PHE:CD1	2.53	0.44
1:B:203:ARG:HA	1:B:234:THR:O	2.17	0.44
1:B:454:LEU:O	1:B:458:ASP:HB2	2.16	0.44
2:D:153:VAL:O	2:D:154:ASP:HB2	2.18	0.44
1:A:393:ARG:NH1	1:A:763:ARG:HD3	2.32	0.44
1:A:367:TYR:O	1:A:371:LYS:HG3	2.18	0.44
1:B:202:PHE:O	1:B:205:ILE:HG12	2.18	0.44
2:D:44:PRO:HG3	2:D:47:ASP:OD2	2.18	0.44
1:B:710:PRO:HA	1:B:713:ARG:HE	1.82	0.44
1:A:695:ARG:HG3	1:A:695:ARG:NH1	2.30	0.44
1:A:450:LEU:HB2	1:A:451:PRO:HD3	2.00	0.44
1:B:652:ASP:N	1:B:653:PRO:CD	2.80	0.44
1:A:273:LEU:O	1:A:275:TYR:N	2.47	0.43
1:B:577:ALA:O	1:B:578:GLN:CB	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:31:ASN:C	2:C:33:LEU:H	2.22	0.43
1:A:317:VAL:HG22	1:A:320:GLY:HA3	1.99	0.43
1:A:296:ARG:HG3	1:A:296:ARG:HH11	1.83	0.43
1:A:562:TRP:CZ2	1:A:596:PHE:HA	2.53	0.43
1:A:425:PHE:O	1:A:429:GLN:HG2	2.18	0.43
1:A:476:LEU:HD22	1:A:550:LEU:CD2	2.48	0.43
1:B:338:THR:O	1:B:342:VAL:HB	2.18	0.43
1:B:445:TYR:CE1	1:B:449:ILE:HG21	2.54	0.43
1:A:387:LEU:H	1:A:387:LEU:CD2	2.31	0.43
2:C:60:GLU:CD	2:C:60:GLU:H	2.21	0.43
1:B:693:PHE:HA	1:B:696:TYR:HB2	2.00	0.43
2:C:153:VAL:HG12	2:C:153:VAL:O	2.18	0.43
1:B:301:LEU:O	1:B:304:VAL:HG13	2.18	0.43
1:A:291:LYS:O	1:A:292:ASN:HB2	2.19	0.43
1:A:453:LEU:HD23	1:A:453:LEU:HA	1.91	0.43
1:A:269:ASN:O	1:A:270:GLU:C	2.57	0.43
1:A:251:SER:N	1:A:271:VAL:HG12	2.32	0.43
1:A:178:THR:CG2	1:A:210:ILE:HD12	2.49	0.43
2:D:116:GLN:HA	2:D:117:PRO:HD2	1.86	0.43
1:A:233:PHE:O	1:A:234:THR:C	2.56	0.43
1:B:246:CYS:O	1:B:266:SER:HB3	2.18	0.43
1:B:664:HIS:HA	1:B:667:HIS:ND1	2.34	0.43
1:A:480:THR:HB	1:A:526:ARG:NH2	2.34	0.43
1:A:411:SER:O	1:A:415:SER:OG	2.33	0.43
1:B:383:ARG:HD3	1:B:417:TRP:HD1	1.84	0.43
1:B:178:THR:HG21	1:B:210:ILE:O	2.18	0.42
1:A:275:TYR:CZ	1:A:296:ARG:HD3	2.53	0.42
1:A:326:LEU:HD22	1:A:360:ILE:HG12	2.01	0.42
1:B:254:SER:O	1:B:255:ASP:HB2	2.19	0.42
1:A:598:GLN:O	1:A:602:GLN:HG3	2.18	0.42
1:A:570:ALA:C	1:A:572:GLU:H	2.21	0.42
1:A:229:SER:O	1:A:230:ASN:HB2	2.20	0.42
1:A:239:SER:O	1:A:241:LEU:HG	2.18	0.42
2:C:47:ASP:O	2:C:48:LYS:CB	2.68	0.42
1:A:174:ALA:C	1:A:176:LYS:H	2.23	0.42
1:A:319:GLY:H	1:A:320:GLY:HA3	1.78	0.42
2:C:49:GLY:HA3	2:C:144:THR:CG2	2.49	0.42
2:C:23:ARG:HH12	2:C:154:ASP:HA	1.85	0.42
2:C:58:PRO:HB2	2:C:60:GLU:OE2	2.20	0.42
1:A:514:PHE:HB2	1:A:544:MET:SD	2.60	0.42
1:B:280:GLN:O	1:B:354:LYS:HE2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:ASP:C	1:A:201:ASP:OD2	2.58	0.42
1:A:274:GLN:HG3	1:A:294:LEU:CD2	2.36	0.42
2:D:21:ASN:CB	2:D:109:ILE:HD13	2.49	0.42
1:B:450:LEU:HA	1:B:450:LEU:HD23	1.77	0.42
1:A:212:PRO:HG2	1:A:215:PHE:HD1	1.85	0.42
1:B:326:LEU:HD22	1:B:360:ILE:HG12	2.01	0.42
1:A:625:PHE:CD2	1:A:696:TYR:CZ	3.07	0.42
1:B:516:ASN:ND2	1:B:518:ASP:N	2.68	0.42
1:A:454:LEU:HD22	1:A:495:LEU:HD12	2.01	0.42
1:A:572:GLU:HA	1:A:572:GLU:OE1	2.20	0.42
2:D:132:ASP:OD2	2:D:135:LYS:CB	2.68	0.42
1:A:777:LEU:HA	1:A:778:PRO:HD3	1.89	0.41
1:B:612:ILE:HB	1:B:613:LEU:HD12	2.02	0.41
1:B:262:LYS:HG3	1:B:284:SER:OG	2.20	0.41
1:B:516:ASN:ND2	1:B:518:ASP:HB2	2.35	0.41
1:B:777:LEU:CD2	1:B:778:PRO:HD2	2.50	0.41
1:B:323:PHE:HA	1:B:331:LEU:O	2.21	0.41
2:D:70:PHE:HE2	2:D:85:VAL:HG21	1.85	0.41
2:D:73:LYS:HB2	2:D:147:TYR:HB3	2.02	0.41
2:D:153:VAL:O	2:D:154:ASP:CB	2.67	0.41
1:B:233:PHE:O	1:B:256:VAL:HG22	2.20	0.41
1:B:306:LEU:HD21	1:B:363:LEU:HD13	2.02	0.41
1:A:169:SER:CB	1:A:170:PRO:HD2	2.48	0.41
1:A:598:GLN:HA	1:A:601:THR:HB	2.03	0.41
1:B:371:LYS:O	1:B:374:ALA:HB3	2.20	0.41
2:C:36:GLN:HE21	2:C:36:GLN:HB2	1.61	0.41
1:B:180:ALA:HB1	1:B:190:LEU:HD23	2.02	0.41
1:B:696:TYR:HB3	1:B:712:LEU:HD13	2.02	0.41
1:B:580:SER:HA	1:B:581:PRO:HD3	1.96	0.41
2:C:105:ILE:O	2:C:109:ILE:HG13	2.20	0.41
1:B:212:PRO:HG2	1:B:215:PHE:CD1	2.56	0.41
2:C:76:HIS:CD2	2:C:77:PRO:CD	3.02	0.41
1:A:248:PHE:O	1:A:249:THR:C	2.59	0.41
1:B:551:LEU:HA	1:B:551:LEU:HD23	1.81	0.41
1:B:212:PRO:HA	1:B:213:PRO:HD3	1.88	0.41
2:D:89:VAL:HG13	2:D:90:ILE:HG23	2.03	0.41
2:C:53:ILE:HD11	2:C:70:PHE:CD1	2.55	0.41
1:B:212:PRO:C	1:B:214:SER:H	2.23	0.41
1:B:297:HIS:HB2	1:B:356:ILE:HB	2.02	0.41
1:A:526:ARG:O	1:A:547:THR:HB	2.21	0.41
1:B:364:PRO:HG2	1:B:367:TYR:HD1	1.79	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:348:HIS:HB3	1:A:355:SER:OG	2.21	0.41
2:D:52:ARG:O	2:D:53:ILE:HG13	2.21	0.41
1:A:624:GLN:NE2	1:A:636:VAL:HG11	2.36	0.41
1:A:730:ILE:HG23	1:A:769:PHE:CE1	2.56	0.41
1:B:516:ASN:HD22	1:B:518:ASP:N	2.19	0.41
1:A:387:LEU:N	1:A:387:LEU:CD2	2.83	0.41
1:A:265:ASN:HA	1:A:287:ASP:O	2.22	0.41
1:B:500:LEU:HD22	1:B:509:ILE:HD11	2.04	0.40
1:A:259:SER:N	1:A:279:THR:HG23	2.36	0.40
2:D:76:HIS:HA	2:D:77:PRO:HD2	1.81	0.40
1:B:366:LYS:HE3	1:B:367:TYR:CE1	2.56	0.40
2:C:87:LEU:HA	2:C:88:PRO:HD3	1.77	0.40
1:A:556:PRO:HG3	1:A:562:TRP:HH2	1.87	0.40
2:D:117:PRO:HG3	2:D:129:TYR:HD1	1.86	0.40
1:A:608:LEU:HD12	1:A:668:ILE:HG12	2.03	0.40
1:B:310:ASP:C	1:B:312:SER:H	2.25	0.40
1:A:709:PRO:HA	1:A:710:PRO:HD3	1.76	0.40
1:A:173:TRP:CH2	1:A:200:ALA:HA	2.56	0.40
1:A:570:ALA:O	1:A:572:GLU:N	2.53	0.40
2:D:66:PRO:O	2:D:90:ILE:HD12	2.21	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	605/621 (97%)	519 (86%)	74 (12%)	12 (2%)	9	46
1	B	612/621 (99%)	537 (88%)	71 (12%)	4 (1%)	26	69
2	C	152/156 (97%)	135 (89%)	16 (10%)	1 (1%)	26	69
2	D	152/156 (97%)	135 (89%)	16 (10%)	1 (1%)	26	69
All	All	1521/1554 (98%)	1326 (87%)	177 (12%)	18 (1%)	16	57

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	274	GLN
1	A	511	ASN
1	A	291	LYS
1	B	511	ASN
1	B	733	SER
2	D	80	ASP
1	A	191	THR
1	A	206	THR
1	A	312	SER
1	A	442	ASP
1	B	200	ALA
1	B	537	GLN
1	A	337	TRP
1	A	510	HIS
1	A	578	GLN
2	C	29	GLU
1	A	571	GLY
1	A	701	ILE

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	484/541 (90%)	448 (93%)	36 (7%)	17	54
1	B	487/541 (90%)	439 (90%)	48 (10%)	10	37
2	C	115/139 (83%)	106 (92%)	9 (8%)	16	51
2	D	115/139 (83%)	104 (90%)	11 (10%)	10	38
All	All	1201/1360 (88%)	1097 (91%)	104 (9%)	13	45

All (104) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	194	ASP
1	A	196	ASP

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Mol	Chain	Res	Type
1	A	270	GLU
1	A	295	ILE
1	A	312	SER
1	A	328	ASP
1	A	335	LEU
1	A	348	HIS
1	A	387	LEU
1	A	395	VAL
1	A	398	SER
1	A	444	TYR
1	A	465	THR
1	A	482	VAL
1	A	486	ASP
1	A	495	LEU
1	A	533	LEU
1	A	536	SER
1	A	552	THR
1	A	557	THR
1	A	567	LEU
1	A	593	LEU
1	A	609	ILE
1	A	630	ASN
1	A	633	HIS
1	A	640	ASP
1	A	650	ILE
1	A	675	THR
1	A	676	ASP
1	A	678	THR
1	A	695	ARG
1	A	696	TYR
1	A	725	GLU
1	A	744	PHE
1	A	753	CYS
1	A	760	SER
1	B	172	ASP
1	B	183	ARG
1	B	189	THR
1	B	190	LEU
1	B	216	MET
1	B	236	SER
1	B	252	THR
1	B	262	LYS

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Mol	Chain	Res	Type
1	B	273	LEU
1	B	294	LEU
1	B	306	LEU
1	B	332	GLN
1	B	336	LEU
1	B	338	THR
1	B	342	VAL
1	B	345	PHE
1	B	369	GLN
1	B	379	VAL
1	B	387	LEU
1	B	388	THR
1	B	392	ILE
1	B	418	LEU
1	B	442	ASP
1	B	465	THR
1	B	484	CYS
1	B	485	THR
1	B	486	ASP
1	B	516	ASN
1	B	518	ASP
1	B	532	LEU
1	B	541	THR
1	B	550	LEU
1	B	557	THR
1	B	567	LEU
1	B	576	THR
1	B	593	LEU
1	B	596	PHE
1	B	604	ARG
1	B	608	LEU
1	B	630	ASN
1	B	670	SER
1	B	675	THR
1	B	707	ASP
1	B	721	GLN
1	B	741	SER
1	B	744	PHE
1	B	771	SER
1	B	775	SER
2	C	7	LEU
2	C	25	ILE

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Mol	Chain	Res	Type
2	C	31	ASN
2	C	60	GLU
2	C	100	LYS
2	C	102	ASP
2	C	104	VAL
2	C	105	ILE
2	C	122	ARG
2	D	1	MET
2	D	36	GLN
2	D	38	LEU
2	D	47	ASP
2	D	67	LYS
2	D	80	ASP
2	D	102	ASP
2	D	104	VAL
2	D	132	ASP
2	D	150	LYS
2	D	151	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	292	ASN
1	A	573	ASN
1	A	631	GLN
1	A	764	HIS
1	B	516	ASN
1	B	664	HIS
1	B	764	HIS
2	C	76	HIS
2	C	78	ASN
2	C	139	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

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	$\# Z > 2$
3	SO4	A	1	-	4,4,4	0.22	0	6,6,6	0.07	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
3	SO4	A	1	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [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	609/621 (98%)	-0.11	7 (1%) 82 76	40, 87, 146, 314	0
1	B	616/621 (99%)	-0.27	0 100 100	45, 79, 124, 187	0
2	C	154/156 (98%)	-0.22	0 100 100	46, 77, 116, 158	0
2	D	154/156 (98%)	-0.19	1 (0%) 90 87	46, 74, 115, 150	0
All	All	1533/1554 (98%)	-0.19	8 (0%) 91 89	40, 81, 136, 314	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	222	ILE	4.2
1	A	221	ASP	2.4
1	A	343	ASP	2.2
1	A	330	TRP	2.2
2	D	147	TYR	2.2
1	A	189	THR	2.1
1	A	613	LEU	2.0
1	A	180	ALA	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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	SO4	A	1	5/5	0.93	0.37	2.58	138,138,138,138	0

6.5 Other polymers [i](#)

There are no such residues in this entry.