



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 04:09 am BST

PDB ID : 6BOG
Title : Crystal structure of RapA, a Swi2/Snf2 protein that recycles RNA polymerase during transcription
Authors : Shaw, G.X.; Gan, J.; Zhou, Y.N.; Zhang, R.; Joachimiak, A.; Jin, D.J.; Ji, X.
Deposited on : 2017-11-20
Resolution : 3.21 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

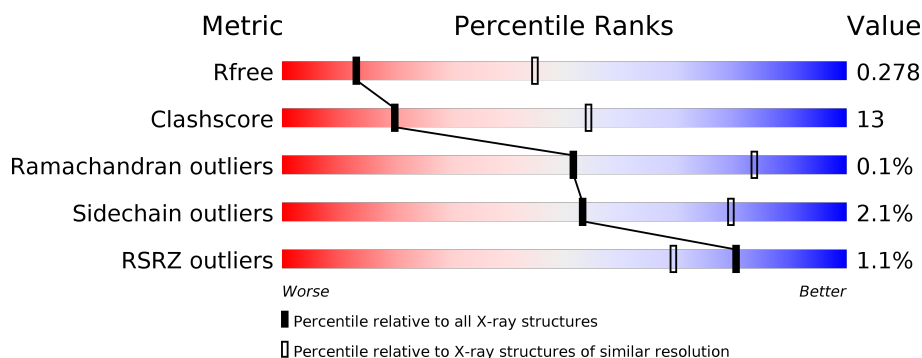
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.21 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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 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	968	<div> <div>%</div> <div> <div></div> <div>70%</div> <div>29%</div> <div>.</div> </div> </div>
1	B	968	<div> <div>%</div> <div> <div></div> <div>73%</div> <div>26%</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
2	SO4	A	1003	-	-	X	-
2	SO4	B	1002	-	-	X	-
2	SO4	B	1003	-	-	X	-

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 15473 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 RNA polymerase-associated protein RapA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	967	Total	C	N	O	S	Se	0	0	0
			7719	4831	1382	1477	6	23			
1	B	967	Total	C	N	O	S	Se	0	0	0
			7719	4831	1382	1477	6	23			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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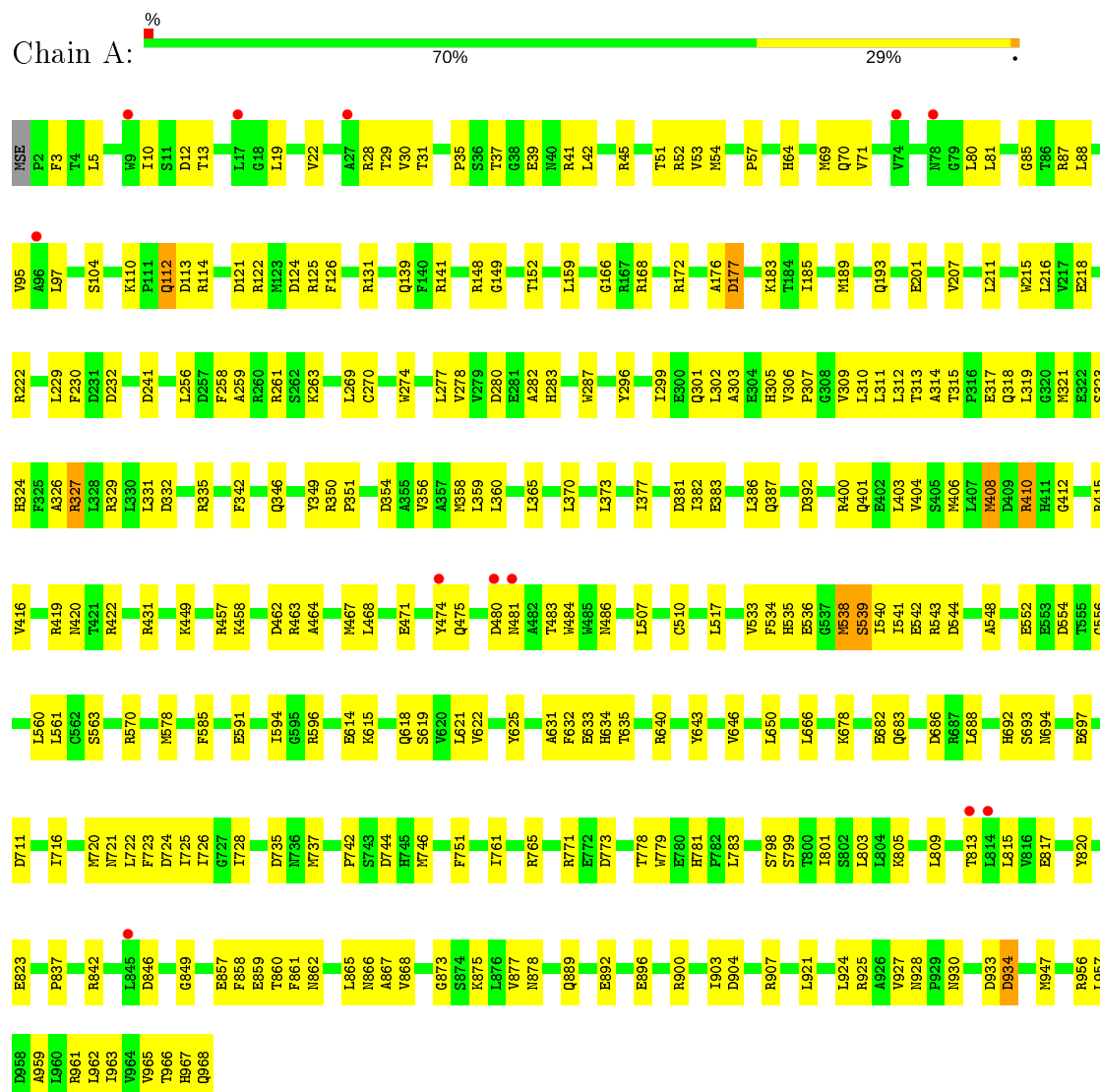
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

3 Residue-property plots

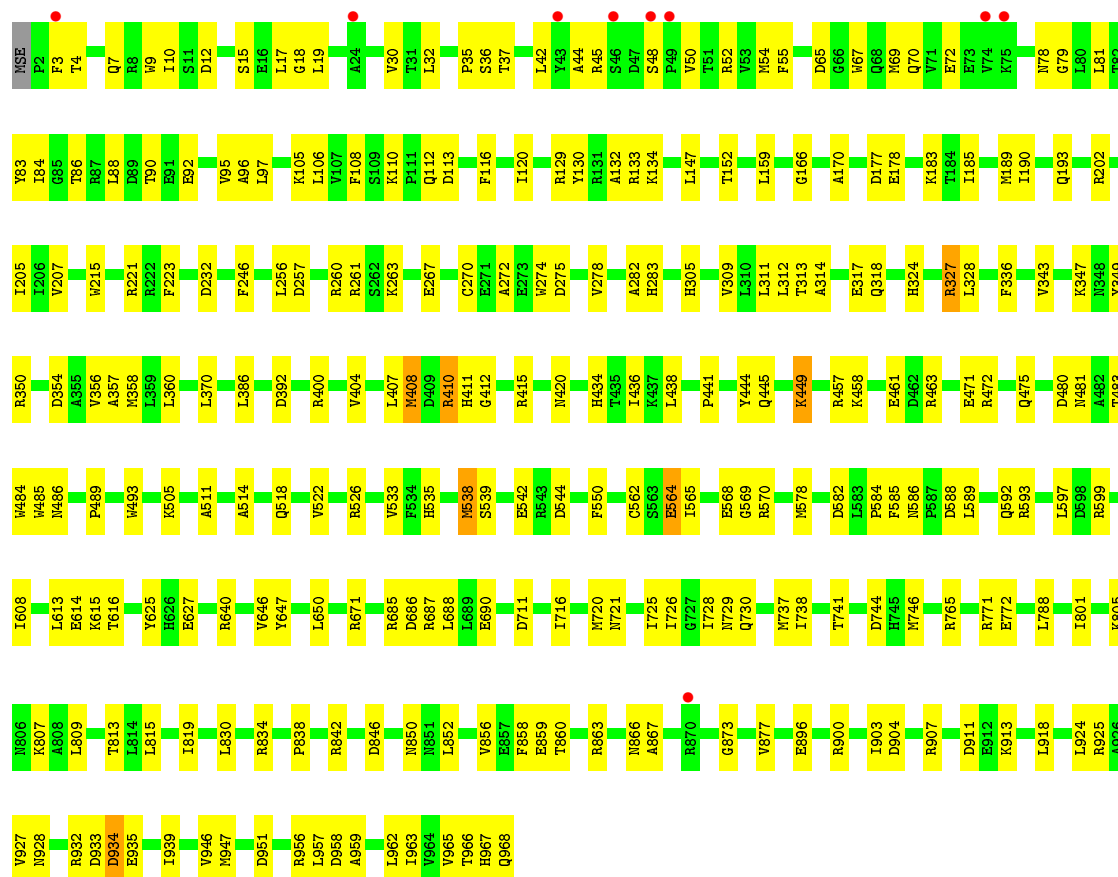
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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



- Molecule 1: RNA polymerase-associated protein RapA





4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	123.86Å 123.86Å 187.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.19 – 3.21 29.19 – 3.20	Depositor EDS
% Data completeness (in resolution range)	98.9 (29.19-3.21) 98.9 (29.19-3.20)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 3.18Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.244 , 0.279 0.245 , 0.278	Depositor DCC
R_{free} test set	994 reflections (2.17%)	wwPDB-VP
Wilson B-factor (Å ²)	91.6	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 38.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.449 for h,-k,-l	Xtriage
Reported twinning fraction	0.500 for h,-k,-l	Depositor
Outliers	0 of 45884 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	15473	wwPDB-VP
Average B, all atoms (Å ²)	105.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.02% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.31	0/7840	0.53	0/10581
1	B	0.29	0/7840	0.53	0/10581
All	All	0.30	0/15680	0.53	0/21162

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

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	7719	0	7583	208	0
1	B	7719	0	7583	183	0
2	A	15	0	0	3	0
2	B	20	0	0	5	0
All	All	15473	0	15166	386	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 13.

All (386) 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:533:VAL:O	1:B:538:MSE:HE1	1.35	1.22
1:B:737:MSE:HE1	1:B:771:ARG:NH1	1.90	0.85
1:A:540:ILE:HD13	1:A:543:ARG:NH2	1.95	0.81
1:B:70:GLN:HB3	1:B:86:THR:HB	1.62	0.80
1:A:540:ILE:CD1	1:A:543:ARG:NH2	2.47	0.77
1:B:67:TRP:HH2	1:B:97:LEU:HD11	1.51	0.74
1:B:7:GLN:NE2	1:B:79:GLY:O	2.20	0.73
1:B:317:GLU:O	1:B:318:GLN:HG2	1.89	0.72
1:A:64:HIS:O	1:B:925:ARG:NH1	2.23	0.72
1:B:627:GLU:O	1:B:671:ARG:NH1	2.23	0.71
1:A:280:ASP:HA	1:A:312:LEU:HB2	1.72	0.71
1:A:536:GLU:O	1:A:536:GLU:HG2	1.89	0.71
1:A:483:THR:OG1	1:A:486:ASN:ND2	2.23	0.71
1:A:114:ARG:NH1	1:A:799:SER:OG	2.24	0.70
1:A:809:LEU:O	1:A:967:HIS:NE2	2.21	0.70
1:B:70:GLN:HB2	1:B:88:LEU:HD21	1.72	0.70
1:A:168:ARG:NH2	1:A:415:ARG:O	2.23	0.70
1:A:110:LYS:HB3	1:A:112:GLN:OE1	1.92	0.69
1:A:166:GLY:HA3	1:A:193:GLN:NE2	2.07	0.69
1:A:538:MSE:HE3	1:A:542:GLU:HB3	1.75	0.69
1:A:925:ARG:NH2	1:A:933:ASP:OD1	2.21	0.69
1:A:172:ARG:HB3	1:A:331:LEU:HD22	1.74	0.68
1:A:722:LEU:HD11	1:A:726:ILE:HD11	1.76	0.68
1:A:817:GLU:OE2	1:A:961:ARG:NH2	2.27	0.68
1:A:131:ARG:NH1	1:A:859:GLU:OE1	2.27	0.68
1:A:646:VAL:HG12	1:A:650:LEU:HB2	1.76	0.67
1:B:410:ARG:HG3	1:B:927:VAL:HG13	1.76	0.67
1:B:221:ARG:HH12	1:B:570:ARG:HH12	1.41	0.67
1:A:324:HIS:HA	1:A:327:ARG:HH11	1.60	0.67
1:B:728:ILE:HG12	1:B:746:MSE:HE1	1.75	0.67
1:B:564:GLU:OE2	1:B:593:ARG:NH2	2.29	0.66
1:A:544:ASP:OD1	1:A:570:ARG:NH1	2.21	0.66
1:A:350:ARG:NH2	1:A:683:GLN:O	2.29	0.66
1:B:511:ALA:N	1:B:582:ASP:OD2	2.29	0.65
1:A:317:GLU:O	1:A:318:GLN:HG2	1.96	0.65
1:A:232:ASP:OD2	1:A:261:ARG:NH1	2.29	0.65
1:B:343:VAL:HG12	1:B:347:LYS:HE2	1.79	0.65
1:A:70:GLN:HB2	1:A:88:LEU:HD21	1.79	0.64
1:A:335:ARG:NE	1:A:930:ASN:OD1	2.28	0.64
1:B:70:GLN:N	1:B:86:THR:O	2.25	0.64
1:B:112:GLN:O	1:B:116:PHE:N	2.25	0.64
1:B:544:ASP:OD1	1:B:570:ARG:NH1	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:588:ASP:OD2	1:B:640:ARG:NE	2.30	0.64
1:B:221:ARG:NH2	1:B:544:ASP:OD2	2.30	0.64
1:A:862:ASN:OD1	1:A:961:ARG:NH1	2.31	0.64
1:B:67:TRP:CH2	1:B:97:LEU:HD11	2.32	0.64
1:A:309:VAL:HG12	1:A:311:LEU:HD13	1.79	0.63
1:A:803:LEU:HD23	1:A:868:VAL:HG12	1.80	0.63
1:A:924:LEU:O	1:A:928:ASN:N	2.32	0.62
1:B:438:LEU:HD13	1:B:489:PRO:HB2	1.81	0.62
1:B:578:MSE:HE3	1:B:597:LEU:HD22	1.80	0.62
1:B:834:ARG:HH21	1:B:913:LYS:NZ	1.97	0.62
1:B:842:ARG:NH1	1:B:858:PHE:HD1	1.98	0.62
1:B:564:GLU:OE2	1:B:593:ARG:NE	2.31	0.61
1:B:584:PRO:HD3	1:B:593:ARG:NH1	2.16	0.61
1:B:685:ARG:HD2	1:B:690:GLU:OE2	2.00	0.61
1:B:813:THR:HG22	1:B:965:VAL:HB	1.83	0.61
1:B:309:VAL:HG12	1:B:311:LEU:HD13	1.82	0.61
1:B:484:TRP:NE1	1:B:518:GLN:OE1	2.31	0.61
1:B:614:GLU:HG2	1:B:615:LYS:HG2	1.82	0.61
1:B:232:ASP:OD1	1:B:261:ARG:NH2	2.34	0.60
1:A:359:LEU:HD21	1:A:400:ARG:HG2	1.82	0.60
1:A:621:LEU:HB3	1:A:625:TYR:CE2	2.36	0.60
1:B:48:SER:HB2	1:B:50:VAL:HG22	1.83	0.60
1:A:646:VAL:HG11	1:A:666:LEU:HD21	1.82	0.60
1:A:323:SER:O	1:A:327:ARG:NH1	2.35	0.60
1:B:129:ARG:NH2	1:B:958:ASP:OD1	2.24	0.60
1:A:35:PRO:HB2	1:A:104:SER:OG	2.02	0.60
1:B:282:ALA:HB3	1:B:313:THR:HB	1.83	0.59
1:B:183:LYS:N	2:B:1001:SO4:O2	2.34	0.59
1:A:538:MSE:HE3	1:A:542:GLU:CB	2.33	0.59
1:B:183:LYS:HB3	1:B:312:LEU:HD23	1.83	0.59
1:A:309:VAL:HG12	1:A:311:LEU:CD1	2.31	0.59
1:A:813:THR:HG22	1:A:965:VAL:HB	1.84	0.59
1:A:966:THR:OG1	1:A:968:GLN:OE1	2.21	0.59
1:B:328:LEU:HB3	1:B:336:PHE:CE2	2.38	0.59
1:A:87:ARG:HB2	1:A:95:VAL:HG21	1.86	0.58
1:A:857:GLU:OE2	1:A:860:THR:HG23	2.04	0.58
1:A:823:GLU:OE2	1:A:956:ARG:NH2	2.36	0.58
1:A:815:LEU:HB2	1:A:963:ILE:HB	1.86	0.58
1:B:270:CYS:HB3	1:B:305:HIS:CG	2.38	0.58
1:A:277:LEU:O	1:A:310:LEU:N	2.33	0.58
1:B:120:ILE:HG21	1:B:801:ILE:HG22	1.84	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:PHE:CD2	1:B:45:ARG:HG2	2.39	0.57
1:A:471:GLU:O	1:A:475:GLN:HG3	2.04	0.57
1:A:517:LEU:HD22	1:A:533:VAL:HG11	1.85	0.57
1:A:185:ILE:HG22	1:A:189:MSE:SE	2.54	0.57
1:B:472:ARG:NH1	1:B:514:ALA:HB3	2.19	0.57
1:B:445:GLN:HB3	1:B:449:LYS:HZ1	1.69	0.57
1:A:110:LYS:NZ	1:A:112:GLN:OE1	2.37	0.57
1:A:335:ARG:NH2	1:A:412:GLY:O	2.38	0.57
1:A:349:TYR:HE1	1:A:688:LEU:HD12	1.70	0.56
1:B:471:GLU:O	1:B:475:GLN:HG3	2.05	0.56
1:B:257:ASP:OD1	1:B:260:ARG:NH1	2.31	0.56
1:A:85:GLY:O	1:A:95:VAL:N	2.38	0.56
1:B:10:ILE:HB	1:B:15:SER:OG	2.04	0.56
1:B:83:TYR:O	1:B:97:LEU:N	2.30	0.56
1:A:422:ARG:NE	1:A:635:THR:HG21	2.19	0.56
1:A:858:PHE:N	2:A:1003:SO4:O3	2.32	0.56
1:A:354:ASP:O	1:A:358:MSE:HG2	2.06	0.56
1:B:415:ARG:HH22	1:B:935:GLU:CD	2.09	0.56
1:B:166:GLY:HA3	1:B:193:GLN:NE2	2.21	0.56
1:B:178:GLU:HB2	1:B:420:ASN:O	2.06	0.55
1:B:925:ARG:NH2	1:B:933:ASP:OD1	2.39	0.55
1:A:431:ARG:NH2	1:A:631:ALA:O	2.40	0.55
1:A:176:ALA:O	1:A:419:ARG:NH1	2.32	0.55
1:B:3:PHE:HD2	1:B:45:ARG:HG2	1.72	0.55
1:B:130:TYR:OH	2:B:1003:SO4:O3	2.25	0.55
1:A:415:ARG:HG2	1:A:415:ARG:HH11	1.71	0.55
1:A:801:ILE:HG13	1:A:962:LEU:HB2	1.89	0.55
1:B:918:LEU:HB2	1:B:939:ILE:HG21	1.89	0.54
1:A:842:ARG:NH1	1:A:858:PHE:HD1	2.06	0.54
1:A:358:MSE:HB2	1:A:365:LEU:HD21	1.90	0.54
1:A:3:PHE:HE2	1:A:45:ARG:HG2	1.72	0.54
1:A:539:SER:OG	1:A:542:GLU:HG3	2.07	0.54
1:A:728:ILE:HG12	1:A:746:MSE:HE1	1.88	0.54
1:B:10:ILE:HG21	1:B:106:LEU:HD11	1.90	0.54
1:B:911:ASP:HB2	1:B:947:MSE:HE2	1.89	0.54
1:A:614:GLU:HG2	1:A:615:LYS:HG2	1.90	0.54
1:B:461:GLU:HG3	1:B:647:TYR:OH	2.07	0.54
1:B:349:TYR:OH	1:B:686:ASP:OD1	2.26	0.54
1:A:19:LEU:HD21	1:A:53:VAL:HG11	1.89	0.54
1:A:857:GLU:HG2	2:A:1003:SO4:O3	2.08	0.53
1:B:538:MSE:HG2	1:B:542:GLU:HB2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:ARG:NE	1:B:272:ALA:O	2.41	0.53
1:B:568:GLU:HA	1:B:599:ARG:NH2	2.22	0.53
1:B:485:TRP:HE1	1:B:526:ARG:NH2	2.06	0.53
1:B:873:GLY:O	1:B:877:VAL:HG23	2.08	0.53
1:A:324:HIS:HA	1:A:327:ARG:NH1	2.24	0.53
1:A:431:ARG:HB2	1:A:633:GLU:HA	1.91	0.53
1:B:807:LYS:HA	1:B:967:HIS:HD2	1.74	0.53
1:B:480:ASP:OD1	1:B:481:ASN:N	2.41	0.53
1:A:585:PHE:O	1:A:625:TYR:OH	2.26	0.53
1:A:370:LEU:HD13	1:A:386:LEU:HD21	1.91	0.52
1:A:771:ARG:NH2	1:A:773:ASP:OD2	2.23	0.52
1:A:907:ARG:HH21	1:A:947:MSE:HB3	1.72	0.52
1:A:535:HIS:O	1:A:538:MSE:HG3	2.09	0.52
1:B:415:ARG:HH11	1:B:415:ARG:HG2	1.74	0.52
1:A:907:ARG:HE	1:A:947:MSE:HG2	1.73	0.52
1:A:507:LEU:HD12	1:A:560:LEU:HB3	1.91	0.52
1:A:900:ARG:HH21	1:A:903:ILE:HG21	1.75	0.52
1:B:54:MSE:HG2	1:B:81:LEU:HD11	1.91	0.52
1:A:726:ILE:HG22	1:A:746:MSE:SE	2.60	0.51
1:A:751:PHE:HE2	1:A:783:LEU:HD22	1.75	0.51
1:B:445:GLN:HB3	1:B:449:LYS:NZ	2.24	0.51
1:B:900:ARG:HH21	1:B:903:ILE:HG12	1.75	0.51
1:A:480:ASP:OD1	1:A:481:ASN:N	2.43	0.51
1:A:172:ARG:HE	1:A:331:LEU:HA	1.76	0.51
1:B:415:ARG:NH2	1:B:935:GLU:OE1	2.43	0.51
1:A:10:ILE:HG23	1:A:53:VAL:HG21	1.92	0.51
1:A:172:ARG:O	1:A:416:VAL:HG23	2.11	0.51
1:B:147:LEU:HG	1:B:159:LEU:HD22	1.93	0.51
1:B:283:HIS:HB3	1:B:314:ALA:O	2.11	0.51
1:B:458:LYS:O	1:B:463:ARG:HD3	2.11	0.51
1:B:408:MSE:HG2	1:B:688:LEU:HD22	1.92	0.51
1:B:350:ARG:NH2	1:B:685:ARG:O	2.43	0.51
1:A:716:ILE:O	1:A:720:MSE:HG3	2.11	0.50
1:B:133:ARG:HH12	1:B:838:PRO:HB2	1.76	0.50
1:B:538:MSE:HG2	1:B:542:GLU:CB	2.41	0.50
1:B:72:GLU:OE2	1:B:86:THR:OG1	2.20	0.50
1:B:815:LEU:HB2	1:B:963:ILE:HB	1.93	0.50
1:A:383:GLU:O	1:A:387:GLN:HG2	2.12	0.50
1:A:552:GLU:HB2	1:A:556:GLY:HA3	1.93	0.50
1:A:110:LYS:HZ2	1:A:110:LYS:HB3	1.75	0.50
1:B:4:THR:HG21	1:B:78:ASN:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:510:CYS:O	1:A:563:SER:HA	2.11	0.50
1:B:805:LYS:HG2	1:B:966:THR:HG23	1.93	0.49
1:B:256:LEU:O	1:B:260:ARG:HB2	2.11	0.49
1:A:408:MSE:SE	1:A:692:HIS:HB2	2.63	0.49
1:B:807:LYS:HA	1:B:967:HIS:CD2	2.48	0.49
1:B:166:GLY:HA2	1:B:190:ILE:HG23	1.95	0.49
1:A:3:PHE:CE2	1:A:45:ARG:HG2	2.47	0.49
1:A:431:ARG:NE	1:A:632:PHE:O	2.42	0.49
1:B:170:ALA:HB3	1:B:932:ARG:HH22	1.78	0.49
1:B:69:MSE:SE	1:B:97:LEU:HD13	2.62	0.49
1:A:207:VAL:HG11	1:A:215:TRP:CD1	2.47	0.49
1:A:678:LYS:O	1:A:682:GLU:HG3	2.13	0.48
1:B:65:ASP:HB2	1:B:67:TRP:NE1	2.29	0.48
1:B:716:ILE:O	1:B:720:MSE:HG2	2.13	0.48
1:B:133:ARG:NH1	1:B:838:PRO:HB2	2.27	0.48
1:B:221:ARG:HD3	1:B:772:GLU:HB3	1.94	0.48
1:B:84:ILE:HA	1:B:95:VAL:O	2.12	0.48
1:A:122:ARG:HB2	1:A:125:ARG:HE	1.78	0.48
1:A:12:ASP:OD2	1:A:51:THR:OG1	2.32	0.48
1:A:39:GLU:OE2	1:A:41:ARG:NH2	2.46	0.48
1:A:805:LYS:HA	1:A:966:THR:O	2.14	0.48
1:A:218:GLU:OE2	1:A:222:ARG:NH2	2.47	0.48
1:A:311:LEU:HD23	1:A:327:ARG:HB3	1.96	0.48
1:A:377:ILE:HD12	1:A:406:MSE:HB3	1.96	0.48
1:B:354:ASP:O	1:B:358:MSE:HG2	2.14	0.48
1:A:585:PHE:CE1	1:A:643:TYR:HE1	2.32	0.48
1:A:866:ASN:OD1	1:A:867:ALA:N	2.47	0.48
1:B:441:PRO:HG2	1:B:444:TYR:CD2	2.49	0.48
1:B:584:PRO:HD3	1:B:593:ARG:CZ	2.44	0.48
1:B:866:ASN:OD1	1:B:867:ALA:N	2.47	0.47
1:B:415:ARG:NH2	1:B:935:GLU:OE2	2.47	0.47
1:B:133:ARG:NH2	1:B:956:ARG:HH22	2.12	0.47
1:B:324:HIS:HA	1:B:327:ARG:NH1	2.28	0.47
1:A:315:THR:HG21	1:A:321:MSE:HE1	1.97	0.47
1:A:781:HIS:CE1	1:A:783:LEU:HG	2.49	0.47
1:B:328:LEU:HB3	1:B:336:PHE:CZ	2.49	0.47
1:A:126:PHE:CE2	1:A:959:ALA:HB1	2.50	0.47
1:A:52:ARG:HG3	1:A:54:MSE:HG3	1.97	0.47
1:A:141:ARG:CZ	1:A:837:PRO:HD2	2.45	0.47
1:A:722:LEU:HD12	1:A:722:LEU:C	2.35	0.46
1:B:966:THR:OG1	1:B:968:GLN:OE1	2.33	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:LEU:HD21	1:A:53:VAL:HG21	1.96	0.46
1:A:159:LEU:HD23	1:A:189:MSE:HE3	1.98	0.46
1:A:400:ARG:O	1:A:404:VAL:HG12	2.15	0.46
1:B:721:ASN:OD1	1:B:725:ILE:HD13	2.16	0.46
1:B:859:GLU:O	1:B:863:ARG:HG3	2.16	0.46
1:B:834:ARG:HH21	1:B:913:LYS:HZ1	1.60	0.46
1:A:13:THR:HG21	1:A:201:GLU:OE2	2.15	0.46
1:A:114:ARG:NH1	1:A:798:SER:OG	2.49	0.46
1:B:132:ALA:HB1	1:B:788:LEU:HD23	1.98	0.46
1:A:152:THR:O	1:A:765:ARG:NH1	2.46	0.46
1:B:434:HIS:N	1:B:608:ILE:O	2.49	0.46
1:A:846:ASP:OD1	1:A:849:GLY:N	2.48	0.46
1:A:934:ASP:OD2	1:B:37:THR:O	2.34	0.46
1:B:370:LEU:HD13	1:B:386:LEU:HD21	1.97	0.46
1:B:69:MSE:HG3	1:B:86:THR:C	2.35	0.46
1:B:96:ALA:O	1:B:97:LEU:HD12	2.16	0.46
1:A:28:ARG:HD2	1:A:241:ASP:OD2	2.16	0.46
1:A:270:CYS:HB3	1:A:305:HIS:CG	2.51	0.46
1:A:69:MSE:SE	1:A:97:LEU:HB2	2.66	0.46
1:B:586:ASN:HD21	1:B:589:LEU:HG	1.81	0.46
1:A:431:ARG:NH2	1:A:634:HIS:O	2.49	0.46
1:A:350:ARG:HB3	1:A:351:PRO:HD3	1.98	0.45
1:A:781:HIS:HE1	1:A:783:LEU:HG	1.80	0.45
1:A:596:ARG:NH2	2:A:1002:SO4:O2	2.44	0.45
1:B:535:HIS:H	1:B:538:MSE:SE	2.49	0.45
1:B:860:THR:HA	1:B:863:ARG:NE	2.31	0.45
1:B:809:LEU:O	1:B:967:HIS:NE2	2.48	0.45
1:B:924:LEU:O	1:B:928:ASN:N	2.46	0.45
1:A:270:CYS:SG	1:A:302:LEU:HA	2.56	0.45
1:A:149:GLY:HA2	1:A:778:THR:HG21	1.98	0.45
1:A:817:GLU:OE1	1:A:842:ARG:NH1	2.50	0.45
1:A:842:ARG:NH1	1:A:858:PHE:HA	2.30	0.45
1:B:404:VAL:HA	1:B:407:LEU:HD12	1.97	0.45
1:B:729:ASN:HB2	1:B:741:THR:OG1	2.16	0.45
1:B:9:TRP:HB2	1:B:32:LEU:HD21	1.98	0.45
1:B:485:TRP:CG	1:B:522:VAL:HG21	2.51	0.45
1:B:900:ARG:HH21	1:B:903:ILE:HG21	1.81	0.45
1:A:177:ASP:HB2	1:A:183:LYS:HG3	1.97	0.45
1:B:202:ARG:O	1:B:275:ASP:N	2.43	0.45
1:A:282:ALA:HB3	1:A:313:THR:HB	1.98	0.45
1:A:744:ASP:N	1:A:744:ASP:OD1	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:809:LEU:HB2	1:B:965:VAL:HG11	1.99	0.45
1:A:342:PHE:O	1:A:346:GLN:HG2	2.17	0.45
1:A:54:MSE:SE	1:A:81:LEU:HD11	2.67	0.45
1:A:861:PHE:O	1:A:865:LEU:HD13	2.16	0.45
1:B:564:GLU:OE2	1:B:593:ARG:CZ	2.65	0.45
1:A:742:PRO:HA	1:A:746:MSE:HE2	1.99	0.45
1:B:730:GLN:HB3	1:B:738:ILE:HG21	1.99	0.45
1:B:830:LEU:HB3	1:B:946:VAL:HG22	1.98	0.45
1:A:873:GLY:O	1:A:877:VAL:HG23	2.17	0.44
1:B:134:LYS:NZ	2:B:1003:SO4:O3	2.43	0.44
1:B:613:LEU:O	1:B:616:THR:HG22	2.15	0.44
1:A:474:TYR:CD2	1:A:484:TRP:HB3	2.52	0.44
1:B:569:GLY:HA3	2:B:1002:SO4:O4	2.18	0.44
1:A:737:MSE:HA	1:A:761:ILE:O	2.17	0.44
1:A:5:LEU:O	1:A:80:LEU:HD21	2.17	0.44
1:A:410:ARG:HG3	1:A:927:VAL:HG13	1.98	0.44
1:B:400:ARG:O	1:B:404:VAL:HG12	2.18	0.44
1:B:729:ASN:N	1:B:741:THR:O	2.49	0.44
1:A:540:ILE:HA	1:A:540:ILE:HD12	1.73	0.44
1:B:15:SER:C	1:B:18:GLY:H	2.19	0.44
1:B:311:LEU:HD23	1:B:327:ARG:HB3	1.99	0.44
1:B:842:ARG:NH1	1:B:858:PHE:HA	2.33	0.44
1:A:112:GLN:HG2	1:A:113:ASP:N	2.33	0.44
1:B:819:ILE:HB	1:B:959:ALA:HB3	1.99	0.44
1:B:842:ARG:NH1	1:B:858:PHE:CD1	2.83	0.44
1:A:274:TRP:O	1:A:307:PRO:HD2	2.18	0.44
1:A:464:ALA:O	1:A:468:LEU:HG	2.17	0.44
1:A:723:PHE:O	1:A:724:ASP:C	2.56	0.44
1:A:382:ILE:O	1:A:386:LEU:N	2.51	0.44
1:A:139:GLN:OE1	1:A:148:ARG:NH1	2.43	0.44
1:A:540:ILE:CD1	1:A:543:ARG:HH21	2.30	0.44
1:A:735:ASP:O	1:A:737:MSE:HG3	2.17	0.44
1:B:170:ALA:O	1:B:932:ARG:NH1	2.39	0.44
1:B:205:ILE:HG23	1:B:278:VAL:HB	1.99	0.44
1:A:415:ARG:NH1	1:A:415:ARG:HG2	2.33	0.43
1:A:269:LEU:HD23	1:A:269:LEU:HA	1.88	0.43
1:A:463:ARG:O	1:A:467:MSE:HG2	2.19	0.43
1:A:401:GLN:NE2	1:A:697:GLU:OE2	2.50	0.43
1:A:721:ASN:OD1	1:A:725:ILE:HD13	2.19	0.43
1:A:296:TYR:HE1	1:A:329:ARG:HH21	1.66	0.43
1:A:842:ARG:HH11	1:A:858:PHE:HD1	1.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:GLN:HG3	1:A:779:TRP:CG	2.54	0.43
1:B:357:ALA:HB1	1:B:687:ARG:NH1	2.34	0.43
1:A:37:THR:C	1:B:934:ASP:OD2	2.57	0.43
1:A:618:GLN:O	1:A:622:VAL:HG23	2.18	0.43
1:B:846:ASP:OD1	1:B:850:ASN:N	2.34	0.43
1:B:90:THR:OG1	1:B:92:GLU:HG3	2.17	0.43
1:A:277:LEU:HB2	1:A:306:VAL:HG11	2.01	0.43
1:A:578:MSE:HE1	1:A:594:ILE:HA	2.01	0.43
1:B:105:LYS:HE2	1:B:105:LYS:HB2	1.87	0.43
1:B:44:ALA:O	1:B:48:SER:HB3	2.17	0.43
1:B:805:LYS:HA	1:B:966:THR:O	2.19	0.43
1:A:37:THR:O	1:B:934:ASP:OD2	2.37	0.43
1:A:640:ARG:HA	1:A:643:TYR:HB3	2.00	0.43
1:B:589:LEU:HA	1:B:592:GLN:HB2	2.01	0.43
1:A:216:LEU:HD11	1:A:229:LEU:HD13	2.00	0.43
1:B:30:VAL:O	1:B:42:LEU:HD12	2.18	0.43
1:B:585:PHE:O	1:B:625:TYR:OH	2.37	0.43
1:A:5:LEU:HD12	1:A:22:VAL:HG12	2.01	0.43
1:A:258:PHE:HA	1:A:261:ARG:HH21	1.82	0.43
1:A:458:LYS:HD2	1:A:462:ASP:HB2	2.00	0.43
1:A:507:LEU:HD23	1:A:578:MSE:HE2	2.01	0.43
1:B:842:ARG:HH12	1:B:858:PHE:HD1	1.64	0.42
1:A:29:THR:HG21	1:A:241:ASP:HB3	2.01	0.42
1:B:350:ARG:HG3	1:B:354:ASP:OD2	2.18	0.42
1:B:646:VAL:HG12	1:B:650:LEU:HB2	2.00	0.42
1:A:370:LEU:HD13	1:A:386:LEU:CD2	2.49	0.42
1:B:410:ARG:HG3	1:B:927:VAL:CG1	2.47	0.42
1:B:562:CYS:SG	1:B:565:ILE:HD13	2.60	0.42
1:A:373:LEU:HD11	1:A:403:LEU:HD11	2.00	0.42
1:B:436:ILE:HG12	1:B:493:TRP:NE1	2.34	0.42
1:B:907:ARG:NH1	1:B:951:ASP:OD1	2.52	0.42
1:A:28:ARG:HB2	1:A:241:ASP:OD2	2.20	0.42
1:A:278:VAL:HG22	1:A:310:LEU:HD12	2.02	0.42
1:A:283:HIS:HB3	1:A:314:ALA:O	2.20	0.42
1:A:30:VAL:O	1:A:42:LEU:HD12	2.20	0.42
1:B:411:HIS:HD1	1:B:412:GLY:H	1.68	0.42
1:A:172:ARG:HH21	1:A:331:LEU:HA	1.85	0.42
1:B:110:LYS:HE2	1:B:113:ASP:OD2	2.20	0.42
1:B:356:VAL:O	1:B:360:LEU:HD12	2.20	0.42
1:A:431:ARG:NH1	1:A:591:GLU:OE2	2.53	0.42
1:A:722:LEU:O	1:A:722:LEU:HD12	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:570:ARG:NH2	2:B:1002:SO4:O2	2.52	0.42
1:B:202:ARG:NH2	1:B:246:PHE:O	2.26	0.42
1:B:726:ILE:HG22	1:B:746:MSE:SE	2.70	0.42
1:A:122:ARG:HB3	1:A:124:ASP:OD1	2.20	0.42
1:A:122:ARG:HB2	1:A:125:ARG:HG3	2.02	0.42
1:A:548:ALA:HA	1:A:771:ARG:HD3	2.02	0.42
1:A:933:ASP:HB2	1:B:35:PRO:O	2.20	0.42
1:A:697:GLU:OE1	1:A:697:GLU:N	2.53	0.41
1:B:147:LEU:HD12	1:B:147:LEU:HA	1.88	0.41
1:B:185:ILE:HG22	1:B:189:MSE:SE	2.69	0.41
1:A:534:PHE:N	1:A:561:LEU:O	2.49	0.41
1:B:272:ALA:HB3	1:B:274:TRP:NE1	2.35	0.41
1:A:31:THR:OG1	1:A:42:LEU:HD13	2.21	0.41
1:A:538:MSE:HB3	1:A:538:MSE:HE3	1.71	0.41
1:A:889:GLN:O	1:A:892:GLU:HB3	2.20	0.41
1:A:309:VAL:CG1	1:A:311:LEU:CD1	2.98	0.41
1:A:382:ILE:HG22	1:A:386:LEU:HB2	2.02	0.41
1:A:349:TYR:CE1	1:A:688:LEU:HD12	2.52	0.41
1:A:540:ILE:CG2	1:A:541:ILE:N	2.81	0.41
1:A:121:ASP:N	1:A:799:SER:O	2.37	0.41
1:A:256:LEU:HD12	1:A:259:ALA:HB3	2.01	0.41
1:A:57:PRO:HA	1:A:71:VAL:HG12	2.01	0.41
1:B:17:LEU:O	1:B:36:SER:OG	2.30	0.41
1:B:263:LYS:O	1:B:267:GLU:N	2.49	0.41
1:B:489:PRO:HD2	1:B:613:LEU:HD21	2.02	0.41
1:B:744:ASP:OD1	1:B:744:ASP:N	2.53	0.41
1:A:875:LYS:HA	1:A:878:ASN:OD1	2.20	0.41
1:B:189:MSE:HG3	1:B:223:PHE:CE1	2.56	0.41
1:B:9:TRP:CZ2	1:B:52:ARG:HD3	2.56	0.41
1:A:306:VAL:HA	1:A:307:PRO:HD3	1.97	0.41
1:A:356:VAL:O	1:A:360:LEU:HD12	2.20	0.41
1:B:207:VAL:HG11	1:B:215:TRP:CD1	2.56	0.41
1:B:627:GLU:O	1:B:671:ARG:HD3	2.21	0.41
1:A:299:ILE:O	1:A:303:ALA:N	2.30	0.41
1:A:263:LYS:HZ1	1:A:301:GLN:HE22	1.68	0.41
1:A:349:TYR:OH	1:A:686:ASP:OD1	2.29	0.41
1:A:319:LEU:HD23	1:A:323:SER:HB3	2.03	0.41
1:A:805:LYS:HG2	1:A:966:THR:HG23	2.03	0.41
1:A:820:TYR:CE2	1:A:957:LEU:HD23	2.56	0.41
1:B:152:THR:O	1:B:765:ARG:NH1	2.52	0.41
1:B:55:PHE:HD2	1:B:83:TYR:CZ	2.39	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:578:MSE:HE1	1:B:593:ARG:O	2.21	0.41
1:B:962:LEU:O	1:B:963:ILE:HD13	2.21	0.41
1:A:332:ASP:HB3	1:A:335:ARG:HB3	2.02	0.40
1:A:349:TYR:OH	1:A:688:LEU:HB2	2.20	0.40
1:A:381:ASP:N	1:A:381:ASP:OD1	2.52	0.40
1:A:207:VAL:HG12	1:A:211:LEU:HB2	2.04	0.40
1:B:505:LYS:HB3	1:B:550:PHE:CZ	2.56	0.40
1:B:586:ASN:ND2	1:B:589:LEU:HG	2.37	0.40
1:B:737:MSE:HE1	1:B:771:ARG:CZ	2.49	0.40
1:B:852:LEU:O	1:B:856:VAL:HG23	2.21	0.40
1:B:957:LEU:HA	1:B:957:LEU:HD22	1.97	0.40
1:B:483:THR:OG1	1:B:486:ASN:ND2	2.54	0.40
1:A:420:ASN:OD1	1:A:693:SER:HB3	2.21	0.40
1:A:615:LYS:HD3	1:A:615:LYS:HA	1.87	0.40
1:A:921:LEU:HD23	1:A:921:LEU:HA	1.83	0.40
1:A:287:TRP:HB3	1:A:326:ALA:HB1	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	965/968 (100%)	930 (96%)	35 (4%)	0	100	100
1	B	965/968 (100%)	928 (96%)	36 (4%)	1 (0%)	51	83
All	All	1930/1936 (100%)	1858 (96%)	71 (4%)	1 (0%)	51	83

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	564	GLU

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	828/805 (103%)	810 (98%)	18 (2%)	52	79
1	B	828/805 (103%)	812 (98%)	16 (2%)	57	81
All	All	1656/1610 (103%)	1622 (98%)	34 (2%)	53	79

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	GLN
1	A	177	ASP
1	A	230	PHE
1	A	327	ARG
1	A	392	ASP
1	A	408	MSE
1	A	410	ARG
1	A	449	LYS
1	A	457	ARG
1	A	538	MSE
1	A	539	SER
1	A	554	ASP
1	A	619	SER
1	A	694	ASN
1	A	711	ASP
1	A	896	GLU
1	A	904	ASP
1	A	934	ASP
1	B	12	ASP
1	B	19	LEU
1	B	108	PHE
1	B	177	ASP
1	B	327	ARG
1	B	392	ASP
1	B	408	MSE
1	B	410	ARG
1	B	449	LYS

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Mol	Chain	Res	Type
1	B	457	ARG
1	B	538	MSE
1	B	539	SER
1	B	711	ASP
1	B	896	GLU
1	B	904	ASP
1	B	934	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	192	HIS
1	A	193	GLN
1	A	694	ASN
1	A	942	ASN
1	B	163	HIS
1	B	475	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

7 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	1002	-	4,4,4	0.15	0	6,6,6	0.13	0
2	SO4	B	1002	-	4,4,4	0.15	0	6,6,6	0.11	0
2	SO4	A	1003	-	4,4,4	0.15	0	6,6,6	0.14	0
2	SO4	B	1001	-	4,4,4	0.14	0	6,6,6	0.13	0
2	SO4	B	1003	-	4,4,4	0.14	0	6,6,6	0.07	0
2	SO4	B	1004	-	4,4,4	0.15	0	6,6,6	0.04	0
2	SO4	A	1001	-	4,4,4	0.14	0	6,6,6	0.18	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.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1002	SO4	1	0
2	B	1002	SO4	2	0
2	A	1003	SO4	2	0
2	B	1001	SO4	1	0
2	B	1003	SO4	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	944/968 (97%)	-0.03	12 (1%) 77 65	58, 105, 147, 170	0
1	B	944/968 (97%)	-0.05	9 (0%) 82 72	43, 99, 151, 199	0
All	All	1888/1936 (97%)	-0.04	21 (1%) 80 69	43, 102, 149, 199	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	49	PRO	5.4
1	A	813	THR	4.6
1	B	3	PHE	4.4
1	B	74	VAL	4.0
1	A	480	ASP	3.9
1	B	46	SER	3.7
1	A	27	ALA	3.6
1	A	74	VAL	3.5
1	B	870	ARG	3.3
1	B	24	ALA	3.1
1	A	17	LEU	3.0
1	B	48	SER	3.0
1	B	75	LYS	2.8
1	A	481	ASN	2.8
1	A	845	LEU	2.7
1	B	43	TYR	2.3
1	A	9	TRP	2.2
1	A	474	TYR	2.1
1	A	78	ASN	2.1
1	A	814	LEU	2.0
1	A	96	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. 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(\AA^2)	Q<0.9
2	SO4	B	1003	5/5	0.90	0.18	113,118,134,134	0
2	SO4	A	1002	5/5	0.91	0.18	74,91,95,99	0
2	SO4	B	1002	5/5	0.94	0.21	74,75,82,103	0
2	SO4	A	1003	5/5	0.95	0.14	95,102,125,133	0
2	SO4	B	1004	5/5	0.97	0.21	84,98,114,122	0
2	SO4	B	1001	5/5	0.98	0.12	64,64,69,78	0
2	SO4	A	1001	5/5	0.99	0.17	56,57,70,88	0

6.5 Other polymers [i](#)

There are no such residues in this entry.