



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 26, 2022 – 06:17 PM JST

PDB ID : 7X20
Title : Crystal structure of non gastric H,K-ATPase alpha2 in (K⁺)E2-AlF state
Authors : Nakanishi, H.; Abe, K.
Deposited on : 2022-02-25
Resolution : 3.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.31.2
buster-report : 1.1.7 (2018)
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) : 2.31.2

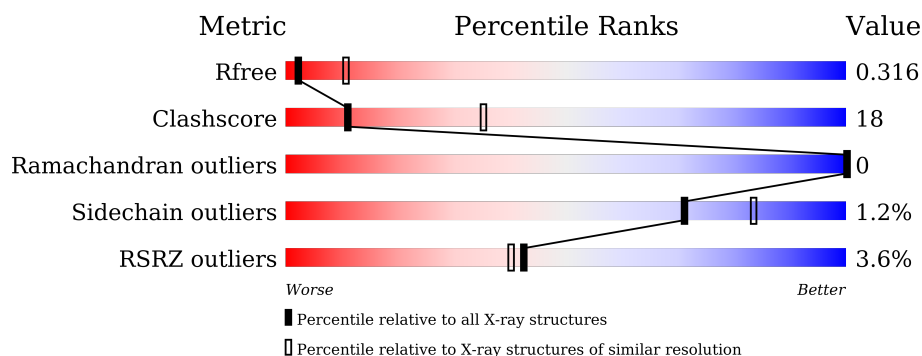
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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

Mol	Chain	Length	Quality of chain
1	A	986	<div> <div>4%</div> <div>62%</div> <div>37%</div> <div>.</div> </div>
2	B	324	<div> <div>2%</div> <div>58%</div> <div>31%</div> <div>11%</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	ALF	A	1101	-	-	X	-
5	NAG	B	401	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10041 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 Potassium-transporting ATPase alpha chain 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	986	Total	C	N	O	S	0	0	0
			7666	4914	1281	1434	37			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	51	GLY	-	expression tag	UNP P54708
A	52	MET	-	expression tag	UNP P54708
A	315	GLY	ASP	conflict	UNP P54708

- Molecule 2 is a protein called Sodium/potassium-transporting ATPase subunit beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	288	Total	C	N	O	S	0	0	0
			2353	1524	385	432	12			

There are 20 discrepancies between the modelled and reference sequences:

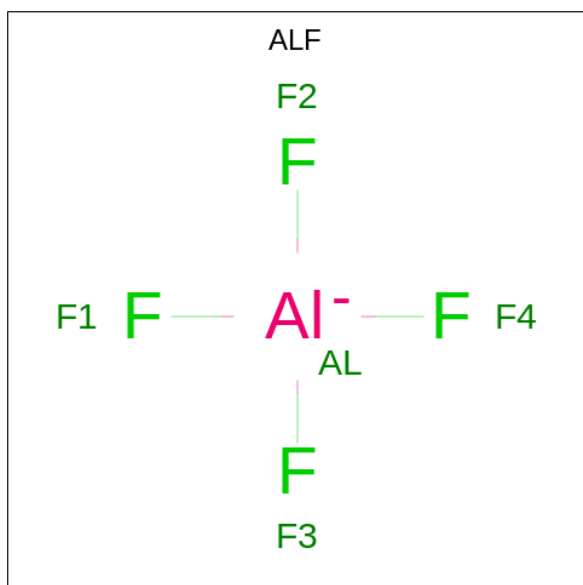
Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	initiating methionine	UNP P07340
B	-18	GLY	-	expression tag	UNP P07340
B	-17	ASP	-	expression tag	UNP P07340
B	-16	TYR	-	expression tag	UNP P07340
B	-15	LYS	-	expression tag	UNP P07340
B	-14	ASP	-	expression tag	UNP P07340
B	-13	ASP	-	expression tag	UNP P07340
B	-12	ASP	-	expression tag	UNP P07340
B	-11	ASP	-	expression tag	UNP P07340
B	-10	LYS	-	expression tag	UNP P07340
B	-9	SER	-	expression tag	UNP P07340
B	-8	SER	-	expression tag	UNP P07340
B	-7	GLY	-	expression tag	UNP P07340

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	GLU	-	expression tag	UNP P07340
B	-5	ASN	-	expression tag	UNP P07340
B	-4	LEU	-	expression tag	UNP P07340
B	-3	TYR	-	expression tag	UNP P07340
B	-2	PHE	-	expression tag	UNP P07340
B	-1	GLN	-	expression tag	UNP P07340
B	0	GLY	-	expression tag	UNP P07340

- Molecule 3 is TETRAFLUOROALUMINATE ION (three-letter code: ALF) (formula: AlF_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Al	F	0	0
			5	1	4		

- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	K	0	0
			2	2		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $\text{C}_8\text{H}_{15}\text{NO}_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			14	8	1	5		

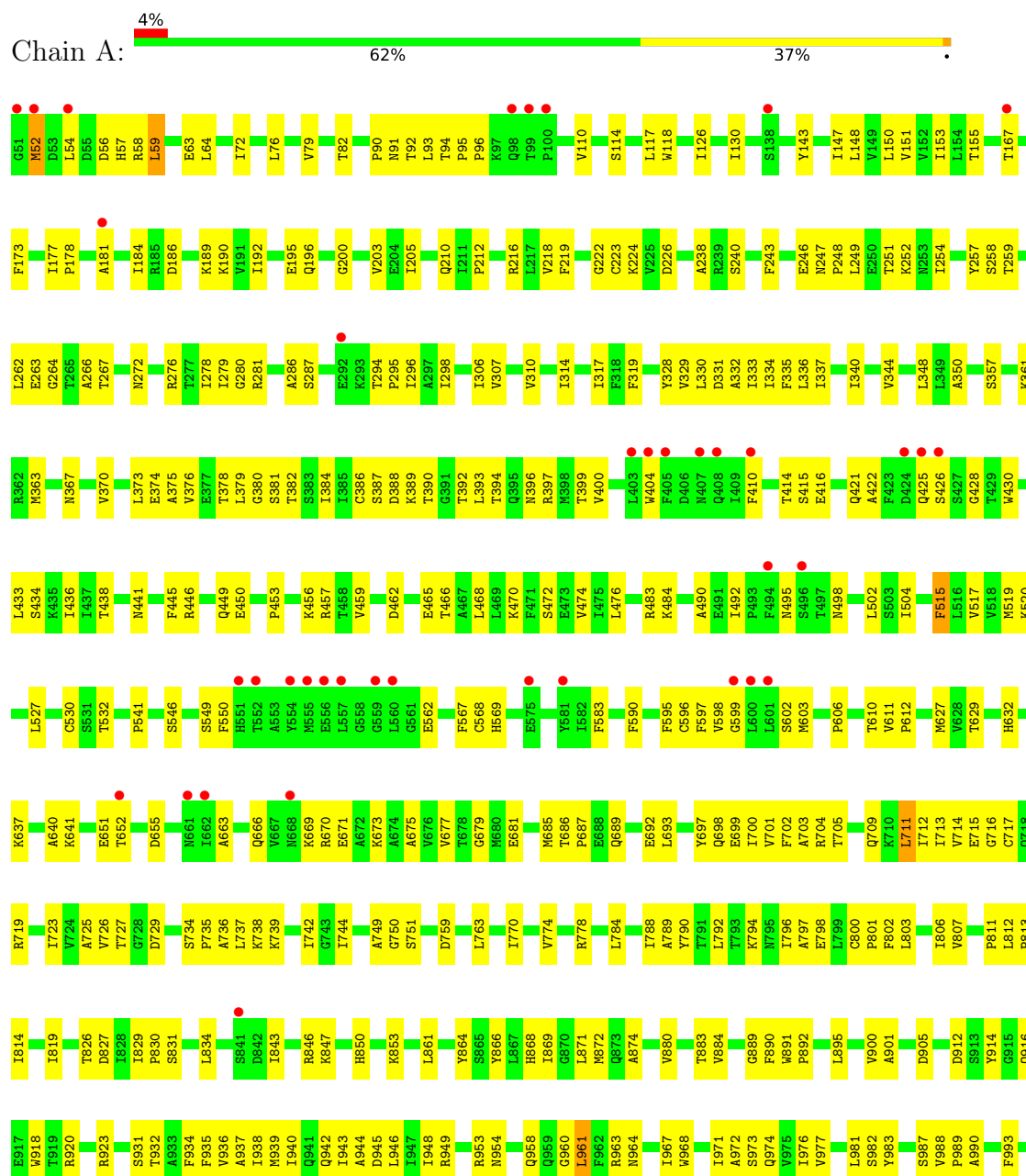
- Molecule 6 is water.

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

3 Residue-property plots

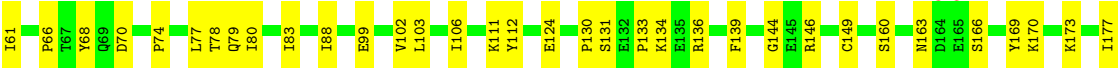
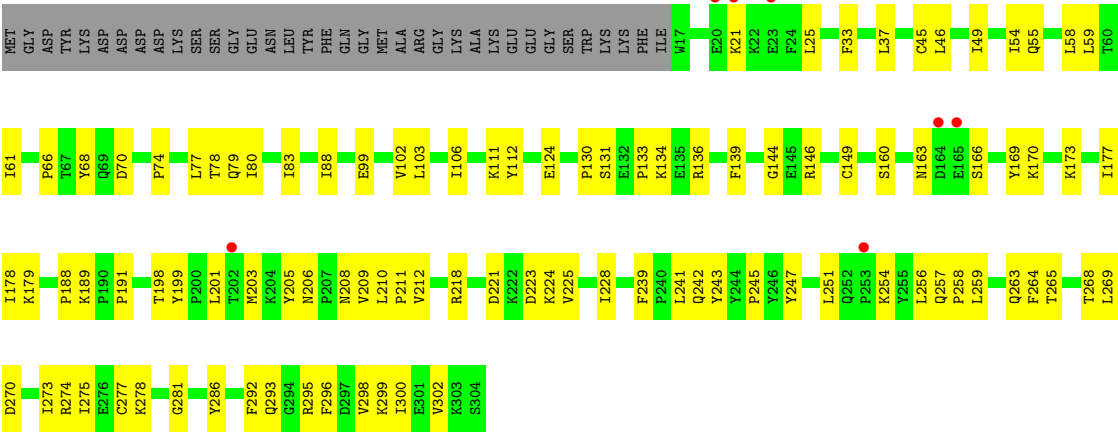
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Potassium-transporting ATPase alpha chain 2





● Molecule 2: Sodium/potassium-transporting ATPase subunit beta-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.97Å 109.06Å 267.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.06 – 3.30 48.06 – 3.13	Depositor EDS
% Data completeness (in resolution range)	72.6 (48.06-3.30) 63.1 (48.06-3.13)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.19_4092	Depositor
R, R_{free}	0.265 , 0.316 0.265 , 0.316	Depositor DCC
R_{free} test set	1276 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	51.1	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 21.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.75	EDS
Total number of atoms	10041	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: K, NAG, ALF

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	0.54	1/7809 (0.0%)	0.77	4/10598 (0.0%)
2	B	0.56	0/2414	0.78	0/3253
All	All	0.55	1/10223 (0.0%)	0.77	4/13851 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	800	CYS	CB-SG	-5.71	1.72	1.81

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	59	LEU	CB-CG-CD1	6.02	121.23	111.00
1	A	711	LEU	CB-CG-CD2	-5.82	101.11	111.00
1	A	59	LEU	CB-CG-CD2	-5.16	102.23	111.00
1	A	379	LEU	CB-CG-CD2	-5.15	102.24	111.00

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	7666	0	7811	291	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2353	0	2335	79	0
3	A	5	0	0	2	0
4	A	2	0	0	0	0
5	B	14	0	13	0	0
6	A	1	0	0	1	0
All	All	10041	0	10159	364	1

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 18.

All (364) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:LEU:HD13	1:A:64:LEU:HB3	1.44	0.97
1:A:226:ASP:HB2	1:A:262:LEU:HD21	1.49	0.94
1:A:686:THR:HG22	1:A:689:GLN:HG2	1.59	0.83
1:A:814:ILE:HG23	1:A:934:PHE:HB3	1.61	0.83
1:A:59:LEU:CD1	1:A:64:LEU:HB3	2.08	0.82
1:A:987:SER:HB2	1:A:993:PHE:HB3	1.59	0.82
1:A:490:ALA:HB3	1:A:504:ILE:HB	1.63	0.81
1:A:872:MET:HG2	2:B:46:LEU:HD11	1.64	0.80
2:B:173:LYS:HB3	2:B:265:THR:HA	1.62	0.80
1:A:943:ILE:HD12	1:A:971:ILE:HG23	1.63	0.78
2:B:88:ILE:HB	2:B:300:ILE:HG22	1.65	0.76
1:A:72:ILE:HD11	1:A:218:VAL:HG21	1.66	0.76
1:A:502:LEU:HD12	1:A:519:MET:HB2	1.68	0.76
1:A:729:ASP:O	1:A:751:SER:N	2.15	0.75
1:A:91:ASN:HB3	1:A:195:GLU:HA	1.68	0.74
1:A:981:LEU:O	1:A:987:SER:HB3	1.87	0.73
1:A:178:PRO:HB3	1:A:210:GLN:HB2	1.71	0.73
1:A:715:GLU:HG2	1:A:739:LYS:HE2	1.69	0.73
2:B:218:ARG:HB2	2:B:221:ASP:HB3	1.71	0.73
1:A:79:VAL:O	1:A:82:THR:OG1	2.05	0.73
2:B:191:PRO:HD3	2:B:281:GLY:HA2	1.72	0.72
1:A:982:SER:O	1:A:987:SER:OG	2.09	0.71
1:A:905:ASP:HB3	1:A:920:ARG:HH11	1.56	0.71
1:A:205:ILE:HG22	1:A:266:ALA:O	1.92	0.70
1:A:450:GLU:N	1:A:450:GLU:OE1	2.25	0.69
1:A:663:ALA:HB3	1:A:666:GLN:HG2	1.75	0.69
1:A:689:GLN:HA	1:A:692:GLU:HB3	1.74	0.68
1:A:436:ILE:HD13	1:A:567:PHE:HB3	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:ASN:OD1	1:A:249:LEU:HB2	1.94	0.67
1:A:278:ILE:HD12	1:A:279:ILE:N	2.09	0.67
1:A:56:ASP:HB2	1:A:59:LEU:HD21	1.77	0.67
1:A:803:LEU:O	1:A:807:VAL:HG22	1.94	0.67
1:A:892:PRO:HG3	2:B:58:LEU:HD11	1.77	0.67
1:A:367:ASN:HA	1:A:763:LEU:HD12	1.78	0.66
1:A:388:ASP:OD2	3:A:1101:ALF:F1	2.04	0.66
1:A:834:LEU:HD11	1:A:949:ARG:HB3	1.78	0.65
1:A:380:GLY:HA2	1:A:774:VAL:HG13	1.79	0.65
1:A:686:THR:HG22	1:A:689:GLN:CG	2.26	0.65
1:A:130:ILE:HG21	1:A:330:LEU:HD13	1.79	0.65
1:A:504:ILE:HG13	1:A:517:VAL:HG23	1.78	0.65
2:B:77:LEU:HB3	2:B:296:PHE:HE1	1.62	0.65
1:A:727:THR:HA	1:A:744:ILE:O	1.97	0.64
1:A:382:THR:HA	1:A:723:ILE:HB	1.78	0.64
1:A:54:LEU:HD12	1:A:54:LEU:O	1.97	0.64
1:A:295:PRO:HA	1:A:298:ILE:HG12	1.80	0.64
1:A:920:ARG:HD2	1:A:923:ARG:HH21	1.63	0.63
1:A:492:ILE:HD12	1:A:502:LEU:HB3	1.79	0.63
1:A:778:ARG:HB3	1:A:843:ILE:HG21	1.80	0.63
1:A:258:SER:OG	1:A:279:ILE:HD13	1.99	0.63
1:A:964:ASN:ND2	1:A:967:ILE:HG23	2.13	0.63
1:A:562:GLU:HB3	1:A:602:SER:HB2	1.80	0.63
1:A:1008:ALA:HA	1:A:1011:ILE:HG22	1.81	0.63
1:A:390:THR:HA	1:A:394:THR:HB	1.81	0.62
2:B:80:ILE:HB	2:B:177:ILE:HB	1.81	0.62
1:A:370:VAL:HG13	1:A:375:ALA:HB3	1.82	0.62
1:A:687:PRO:HB3	1:A:719:ARG:HH12	1.64	0.62
1:A:997:ARG:HH21	2:B:68:TYR:HD2	1.48	0.62
1:A:294:THR:O	1:A:298:ILE:HG23	2.00	0.61
1:A:827:ASP:O	1:A:831:SER:HB3	2.00	0.61
1:A:744:ILE:HD13	1:A:770:ILE:HD12	1.82	0.61
1:A:357:SER:HB3	1:A:373:LEU:HD21	1.81	0.61
1:A:224:LYS:HA	1:A:238:ALA:HA	1.81	0.61
1:A:671:GLU:OE1	1:A:671:GLU:N	2.32	0.61
2:B:239:PHE:HD1	2:B:258:PRO:HB2	1.65	0.61
2:B:144:GLY:O	2:B:146:ARG:NH1	2.33	0.60
1:A:400:VAL:HG12	1:A:603:MET:HG2	1.83	0.60
1:A:797:ALA:HA	1:A:874:ALA:HB2	1.82	0.60
2:B:160:SER:HA	2:B:166:SER:OG	2.01	0.60
1:A:143:TYR:O	1:A:147:ILE:HG13	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:627:MET:HB3	1:A:701:VAL:HG12	1.84	0.60
1:A:711:LEU:HA	1:A:714:VAL:HG12	1.84	0.60
1:A:742:ILE:HG13	1:A:759:ASP:HB2	1.84	0.60
2:B:112:TYR:CZ	2:B:259:LEU:HD11	2.37	0.60
1:A:155:THR:HG22	1:A:350:ALA:HB2	1.84	0.59
1:A:889:GLY:HA2	1:A:914:TYR:CE2	2.36	0.59
1:A:932:THR:O	1:A:936:VAL:HG23	2.03	0.59
1:A:850:HIS:HB3	1:A:853:LYS:HB2	1.83	0.59
2:B:33:PHE:CZ	2:B:37:LEU:HD12	2.37	0.59
1:A:296:ILE:HG21	1:A:374:GLU:HB3	1.84	0.58
1:A:1027:GLY:H	1:A:1031:ASP:CB	2.16	0.58
1:A:453:PRO:O	1:A:457:ARG:HG3	2.03	0.58
1:A:770:ILE:O	1:A:774:VAL:HG23	2.03	0.58
2:B:224:LYS:HB3	2:B:268:THR:OG1	2.04	0.58
1:A:798:GLU:HB3	1:A:819:ILE:HD12	1.85	0.58
1:A:397:ARG:O	1:A:399:THR:HG23	2.03	0.57
1:A:960:GLY:HA3	1:A:963:ARG:HG2	1.85	0.57
1:A:726:VAL:HG23	1:A:737:LEU:HD23	1.85	0.57
1:A:203:VAL:HG11	1:A:212:PRO:HG2	1.85	0.57
1:A:880:VAL:O	1:A:884:VAL:HG12	2.03	0.57
1:A:59:LEU:HD13	1:A:64:LEU:CB	2.27	0.57
1:A:517:VAL:HG13	1:A:595:PHE:HE2	1.70	0.57
1:A:734:SER:OG	1:A:735:PRO:HD3	2.05	0.57
1:A:441:ASN:HB2	1:A:483:ARG:NH2	2.20	0.56
1:A:632:HIS:CD2	1:A:704:ARG:HD2	2.40	0.56
1:A:774:VAL:HG12	1:A:778:ARG:HD2	1.87	0.56
2:B:77:LEU:HB3	2:B:296:PHE:CE1	2.40	0.56
1:A:679:GLY:N	1:A:703:ALA:O	2.36	0.56
1:A:96:PRO:HB3	1:A:287:SER:HB2	1.88	0.56
1:A:669:LYS:HB3	1:A:697:TYR:OH	2.06	0.56
2:B:55:GLN:O	2:B:59:LEU:HG	2.04	0.56
1:A:796:ILE:HG21	1:A:871:LEU:HD12	1.86	0.56
1:A:891:TRP:HB3	1:A:892:PRO:HD2	1.87	0.55
1:A:410:PHE:HE2	1:A:425:GLN:HG3	1.70	0.55
2:B:79:GLN:HE22	2:B:292:PHE:HE2	1.54	0.55
1:A:441:ASN:HD21	1:A:468:LEU:HB2	1.71	0.55
1:A:532:THR:O	1:A:596:CYS:HA	2.06	0.55
1:A:725:ALA:HA	1:A:742:ILE:O	2.07	0.55
1:A:334:ILE:HD12	1:A:337:ILE:HD12	1.88	0.55
1:A:416:GLU:HG2	1:A:456:LYS:HE3	1.89	0.55
1:A:433:LEU:HD12	1:A:599:GLY:HA3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:629:THR:O	1:A:705:THR:HG22	2.07	0.55
2:B:160:SER:HB2	2:B:169:TYR:HE1	1.72	0.54
1:A:415:SER:OG	1:A:416:GLU:N	2.40	0.54
1:A:889:GLY:HA2	1:A:914:TYR:HE2	1.73	0.54
1:A:331:ASP:O	1:A:334:ILE:HG22	2.07	0.54
2:B:33:PHE:CE1	2:B:37:LEU:HD12	2.43	0.54
2:B:179:LYS:HD2	2:B:257:GLN:NE2	2.23	0.53
1:A:310:VAL:HG21	1:A:866:TYR:HE2	1.73	0.53
1:A:726:VAL:HG21	1:A:736:ALA:O	2.08	0.53
1:A:248:PRO:O	1:A:254:ILE:HD11	2.08	0.53
1:A:306:ILE:HG13	1:A:307:VAL:N	2.23	0.53
1:A:651:GLU:HB3	1:A:655:ASP:HB2	1.90	0.53
2:B:254:LYS:O	2:B:256:LEU:HD22	2.09	0.53
2:B:274:ARG:HG2	2:B:299:LYS:HG2	1.89	0.53
1:A:381:SER:O	1:A:723:ILE:HD13	2.08	0.53
1:A:801:PRO:HB3	1:A:813:PRO:HB2	1.91	0.53
1:A:1027:GLY:H	1:A:1031:ASP:HB3	1.72	0.53
1:A:184:ILE:HG12	1:A:189:LYS:HD3	1.91	0.53
1:A:386:CYS:HB2	1:A:726:VAL:HG12	1.90	0.53
1:A:961:LEU:HD12	1:A:968:TRP:HZ2	1.74	0.53
1:A:446:ARG:O	1:A:449:GLN:HG3	2.10	0.52
1:A:964:ASN:HB3	1:A:967:ILE:HG12	1.91	0.52
1:A:332:ALA:HA	1:A:335:PHE:HD2	1.74	0.52
1:A:709:GLN:HA	1:A:712:ILE:HB	1.92	0.52
1:A:990:ALA:O	1:A:993:PHE:HB2	2.10	0.52
1:A:465:GLU:HG3	1:A:520:LYS:NZ	2.25	0.52
1:A:517:VAL:HG13	1:A:595:PHE:CE2	2.45	0.52
1:A:546:SER:HA	1:A:549:SER:HB2	1.92	0.51
2:B:112:TYR:CZ	2:B:256:LEU:HD12	2.45	0.51
1:A:685:MET:HG3	1:A:689:GLN:HG3	1.92	0.51
1:A:59:LEU:CD1	1:A:64:LEU:HD23	2.40	0.51
1:A:389:LYS:O	1:A:393:LEU:N	2.42	0.51
1:A:723:ILE:H	1:A:723:ILE:HD12	1.76	0.51
2:B:298:VAL:HG12	2:B:300:ILE:HG23	1.92	0.51
1:A:794:LYS:HE3	6:A:1201:HOH:O	2.11	0.51
2:B:205:TYR:HE2	2:B:208:ASN:HB2	1.76	0.51
1:A:310:VAL:O	1:A:314:ILE:HG12	2.11	0.51
1:A:954:ASN:ND2	1:A:958:GLN:HG3	2.26	0.51
1:A:697:TYR:O	1:A:700:ILE:HD11	2.10	0.51
1:A:988:VAL:HB	1:A:989:PRO:CD	2.41	0.51
1:A:880:VAL:O	1:A:883:THR:HG22	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:74:PRO:HB2	2:B:293:GLN:HG3	1.93	0.50
1:A:222:GLY:O	1:A:264:GLY:HA3	2.11	0.50
2:B:218:ARG:O	2:B:221:ASP:HB3	2.12	0.50
1:A:380:GLY:O	1:A:381:SER:OG	2.29	0.50
1:A:729:ASP:HB2	1:A:750:GLY:HA2	1.93	0.50
1:A:677:VAL:HG13	1:A:681:GLU:HG3	1.94	0.50
1:A:686:THR:CG2	1:A:689:GLN:HG2	2.36	0.50
1:A:569:HIS:O	1:A:596:CYS:HB3	2.11	0.50
1:A:944:ALA:O	1:A:948:ILE:HG23	2.11	0.50
2:B:209:VAL:O	2:B:211:PRO:HD3	2.11	0.50
1:A:57:HIS:CE1	1:A:246:GLU:O	2.65	0.50
1:A:421:GLN:HG2	1:A:422:ALA:N	2.27	0.50
1:A:652:THR:HG23	1:A:655:ASP:H	1.77	0.50
1:A:389:LYS:HG3	1:A:390:THR:N	2.27	0.49
1:A:843:ILE:HG23	1:A:846:ARG:HH21	1.77	0.49
1:A:117:LEU:CD1	1:A:344:VAL:HG11	2.42	0.49
1:A:240:SER:N	1:A:252:LYS:O	2.46	0.49
1:A:778:ARG:HE	1:A:843:ILE:HG22	1.77	0.49
2:B:37:LEU:HD23	2:B:37:LEU:O	2.13	0.49
1:A:173:PHE:O	1:A:177:ILE:HG12	2.13	0.49
1:A:262:LEU:N	1:A:262:LEU:HD22	2.28	0.49
1:A:438:THR:O	1:A:483:ARG:NE	2.46	0.49
2:B:277:CYS:H	2:B:296:PHE:HB2	1.77	0.49
1:A:890:PHE:HE2	1:A:918:TRP:CE2	2.31	0.49
2:B:78:THR:HG23	2:B:179:LYS:HE2	1.93	0.49
2:B:160:SER:HB2	2:B:169:TYR:CE1	2.48	0.49
1:A:223:CYS:HA	1:A:263:GLU:O	2.12	0.48
1:A:459:VAL:HG21	1:A:470:LYS:HE3	1.94	0.48
1:A:716:GLY:O	1:A:719:ARG:HG2	2.14	0.48
1:A:802:PHE:O	1:A:806:ILE:HG22	2.13	0.48
1:A:569:HIS:ND1	1:A:598:VAL:HG22	2.28	0.48
2:B:83:ILE:HG21	2:B:88:ILE:HG12	1.95	0.48
2:B:131:SER:HB3	2:B:242:GLN:HB3	1.94	0.48
1:A:388:ASP:O	1:A:392:THR:OG1	2.28	0.48
1:A:390:THR:O	3:A:1101:ALF:F1	2.21	0.48
1:A:226:ASP:O	1:A:259:THR:HB	2.14	0.48
1:A:248:PRO:HA	1:A:251:THR:OG1	2.13	0.48
1:A:789:ALA:HB1	1:A:949:ARG:HH21	1.79	0.48
1:A:126:ILE:O	1:A:130:ILE:HG12	2.13	0.48
1:A:900:VAL:HG13	1:A:901:ALA:H	1.79	0.48
2:B:45:CYS:O	2:B:49:ILE:HG13	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:410:PHE:CE2	1:A:425:GLN:HG3	2.49	0.47
1:A:789:ALA:HB2	1:A:861:LEU:HD21	1.97	0.47
1:A:939:MET:HA	1:A:942:GLN:HB2	1.97	0.47
1:A:997:ARG:HH22	2:B:70:ASP:HB2	1.80	0.47
1:A:430:TRP:HA	1:A:433:LEU:HB3	1.97	0.47
1:A:438:THR:HG22	1:A:472:SER:HB2	1.95	0.47
1:A:52:MET:HG2	1:A:272:ASN:HD21	1.80	0.47
1:A:1027:GLY:O	1:A:1032:LYS:HE3	2.15	0.47
2:B:264:PHE:CZ	2:B:275:ILE:HD12	2.50	0.47
1:A:257:TYR:HB3	1:A:279:ILE:HG12	1.96	0.47
1:A:811:PRO:HD2	1:A:895:LEU:HD22	1.97	0.47
1:A:276:ARG:O	1:A:281:ARG:NH2	2.47	0.47
2:B:99:GLU:O	2:B:103:LEU:HD23	2.14	0.47
2:B:112:TYR:CE1	2:B:256:LEU:HD12	2.50	0.47
1:A:880:VAL:HG21	1:A:937:ALA:HB2	1.96	0.47
1:A:387:SER:OG	1:A:393:LEU:HD13	2.14	0.46
2:B:112:TYR:CE1	2:B:259:LEU:HD11	2.50	0.46
2:B:201:LEU:O	2:B:203:MET:HG2	2.15	0.46
1:A:76:LEU:HD22	1:A:200:GLY:O	2.15	0.46
1:A:286:ALA:O	1:A:738:LYS:HG3	2.15	0.46
1:A:936:VAL:O	1:A:940:ILE:HG12	2.15	0.46
1:A:258:SER:HG	1:A:279:ILE:HD13	1.80	0.46
1:A:788:ILE:O	1:A:792:LEU:HG	2.15	0.46
1:A:829:ILE:HD13	1:A:973:SER:HB2	1.97	0.46
1:A:110:VAL:HG12	1:A:110:VAL:O	2.16	0.46
1:A:834:LEU:HA	1:A:834:LEU:HD23	1.65	0.46
1:A:987:SER:CB	1:A:993:PHE:HB3	2.39	0.46
1:A:310:VAL:HG21	1:A:866:TYR:CE2	2.51	0.46
1:A:778:ARG:CB	1:A:843:ILE:HG21	2.44	0.46
2:B:54:ILE:HD12	2:B:54:ILE:HG23	1.67	0.46
1:A:348:LEU:HD21	1:A:788:ILE:HG12	1.98	0.46
1:A:296:ILE:CG2	1:A:374:GLU:HB3	2.45	0.46
2:B:178:ILE:HG21	2:B:212:VAL:HG11	1.97	0.46
1:A:118:TRP:CE3	1:A:118:TRP:HA	2.51	0.45
1:A:192:ILE:HD12	1:A:196:GLN:HB3	1.98	0.45
1:A:370:VAL:CG1	1:A:375:ALA:HB3	2.46	0.45
1:A:796:ILE:HD13	1:A:796:ILE:HA	1.74	0.45
1:A:1008:ALA:HA	1:A:1011:ILE:CG2	2.44	0.45
2:B:179:LYS:HD2	2:B:257:GLN:HE22	1.81	0.45
1:A:330:LEU:O	1:A:333:ILE:HG13	2.16	0.45
1:A:972:ALA:O	1:A:976:ILE:HG12	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:58:LEU:O	2:B:61:ILE:HG13	2.16	0.45
1:A:363:MET:HE2	1:A:376:VAL:HG12	1.98	0.45
1:A:370:VAL:HG11	1:A:376:VAL:HG13	1.97	0.45
1:A:492:ILE:CD1	1:A:502:LEU:HB3	2.46	0.45
1:A:912:ASP:OD1	1:A:916:GLN:HB2	2.16	0.45
1:A:973:SER:HA	1:A:976:ILE:HD11	1.97	0.45
2:B:239:PHE:CD1	2:B:258:PRO:HB2	2.48	0.45
1:A:380:GLY:C	1:A:382:THR:H	2.20	0.45
1:A:826:THR:HA	1:A:977:VAL:HG11	1.98	0.45
1:A:434:SER:HB2	1:A:476:LEU:HD11	1.98	0.45
2:B:189:LYS:HA	2:B:189:LYS:HD2	1.69	0.45
2:B:218:ARG:HB2	2:B:221:ASP:CB	2.43	0.45
1:A:90:PRO:HB2	1:A:92:THR:HG22	1.99	0.45
1:A:784:LEU:HD23	1:A:784:LEU:HA	1.72	0.45
1:A:150:LEU:O	1:A:153:ILE:HG22	2.16	0.45
1:A:317:ILE:HD13	1:A:317:ILE:HA	1.72	0.45
2:B:134:LYS:O	2:B:245:PRO:HD3	2.17	0.45
1:A:147:ILE:O	1:A:151:VAL:HG13	2.17	0.44
1:A:864:TYR:O	1:A:868:HIS:HB2	2.18	0.44
1:A:942:GLN:HB3	1:A:974:GLN:HE22	1.82	0.44
1:A:114:SER:O	1:A:118:TRP:CD1	2.70	0.44
1:A:404:TRP:CH2	1:A:550:PHE:HA	2.52	0.44
1:A:462:ASP:O	1:A:466:THR:HG23	2.16	0.44
1:A:568:CYS:HA	1:A:598:VAL:HG23	1.98	0.44
1:A:190:LYS:HD3	1:A:190:LYS:HA	1.76	0.44
1:A:813:PRO:O	1:A:931:SER:HA	2.17	0.44
1:A:1024:LEU:O	1:A:1026:PRO:HD3	2.18	0.44
2:B:133:PRO:HG3	2:B:241:LEU:HB3	1.99	0.44
2:B:136:ARG:NH2	2:B:149:CYS:SG	2.91	0.44
1:A:953:ARG:HH12	1:A:1023:ARG:HA	1.80	0.44
1:A:76:LEU:HA	1:A:76:LEU:HD12	1.60	0.44
1:A:117:LEU:HD11	1:A:344:VAL:HG11	2.00	0.44
1:A:181:ALA:HB2	1:A:212:PRO:HB3	2.00	0.44
1:A:184:ILE:HD11	1:A:189:LYS:HE2	1.99	0.44
1:A:530:CYS:SG	1:A:597:PHE:N	2.91	0.44
2:B:170:LYS:HE2	2:B:263:GLN:HG3	1.99	0.44
1:A:57:HIS:ND1	1:A:243:PHE:HE1	2.16	0.44
1:A:1011:ILE:HG23	1:A:1012:TRP:H	1.83	0.44
1:A:414:THR:HB	1:A:606:PRO:CB	2.48	0.44
1:A:675:ALA:HB3	1:A:700:ILE:HG23	1.99	0.44
2:B:102:VAL:O	2:B:106:ILE:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:223:ASP:C	2:B:225:VAL:H	2.21	0.43
1:A:386:CYS:O	1:A:726:VAL:HA	2.17	0.43
1:A:802:PHE:CD1	1:A:812:LEU:HD11	2.52	0.43
1:A:1011:ILE:HG23	1:A:1012:TRP:N	2.33	0.43
1:A:997:ARG:HE	2:B:68:TYR:HE2	1.61	0.43
1:A:219:PHE:HB2	1:A:267:THR:HB	2.00	0.43
2:B:170:LYS:NZ	2:B:263:GLN:HB2	2.33	0.43
1:A:319:PHE:HD2	1:A:336:LEU:HD13	1.82	0.43
1:A:935:PHE:O	1:A:938:ILE:HG13	2.18	0.43
2:B:269:LEU:HD12	2:B:302:VAL:HB	2.01	0.43
1:A:702:PHE:CD2	1:A:713:ILE:HD12	2.54	0.43
1:A:530:CYS:SG	1:A:597:PHE:HB2	2.59	0.43
2:B:130:PRO:HB2	2:B:205:TYR:CE2	2.53	0.43
2:B:170:LYS:HE2	2:B:263:GLN:CG	2.49	0.43
1:A:94:THR:HA	1:A:95:PRO:HD3	1.88	0.43
1:A:673:LYS:HE3	1:A:698:GLN:NE2	2.34	0.43
1:A:702:PHE:CE1	1:A:713:ILE:HD11	2.53	0.43
1:A:216:ARG:HA	1:A:254:ILE:HG22	2.01	0.43
1:A:361:LYS:HD3	1:A:361:LYS:HA	1.75	0.43
1:A:905:ASP:HB3	1:A:920:ARG:NH1	2.29	0.42
2:B:124:GLU:CD	2:B:134:LYS:HZ3	2.22	0.42
1:A:192:ILE:HD12	1:A:196:GLN:OE1	2.19	0.42
1:A:433:LEU:HA	1:A:436:ILE:HG22	2.01	0.42
1:A:698:GLN:HG2	1:A:699:GLU:HG3	1.99	0.42
1:A:830:PRO:HB3	1:A:946:LEU:HD22	2.02	0.42
2:B:54:ILE:HD13	2:B:54:ILE:HA	1.87	0.42
1:A:332:ALA:O	1:A:335:PHE:HB2	2.19	0.42
1:A:340:ILE:HD13	1:A:340:ILE:HG21	1.74	0.42
2:B:111:LYS:HG2	2:B:256:LEU:HD11	2.01	0.42
1:A:459:VAL:HG11	1:A:466:THR:HG22	2.01	0.42
1:A:52:MET:HB2	1:A:54:LEU:HG	2.02	0.42
1:A:389:LYS:NZ	1:A:390:THR:HG23	2.34	0.42
1:A:400:VAL:HG12	1:A:603:MET:CG	2.49	0.42
1:A:295:PRO:CG	1:A:378:THR:HG23	2.49	0.42
1:A:449:GLN:N	1:A:450:GLU:OE1	2.52	0.42
2:B:243:TYR:CD1	2:B:258:PRO:HG3	2.55	0.42
1:A:790:TYR:CE1	1:A:942:GLN:NE2	2.88	0.42
1:A:93:LEU:HD23	1:A:280:GLY:HA2	2.01	0.42
2:B:163:ASN:N	2:B:166:SER:HB2	2.34	0.42
2:B:273:ILE:O	2:B:299:LYS:HA	2.20	0.42
1:A:329:VAL:O	1:A:333:ILE:HG23	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:610:THR:C	1:A:612:PRO:HD2	2.40	0.42
1:A:426:SER:C	1:A:428:GLY:H	2.23	0.42
1:A:453:PRO:HG2	1:A:456:LYS:HD2	2.02	0.42
1:A:257:TYR:CG	1:A:279:ILE:HD11	2.55	0.41
1:A:393:LEU:HG	1:A:611:VAL:HG21	2.01	0.41
1:A:627:MET:HE2	1:A:640:ALA:HB2	2.01	0.41
2:B:80:ILE:HD13	2:B:80:ILE:HA	1.84	0.41
2:B:228:ILE:HG22	2:B:264:PHE:HD1	1.85	0.41
1:A:515:PHE:HD1	1:A:515:PHE:HA	1.74	0.41
1:A:814:ILE:CG2	1:A:934:PHE:HB3	2.41	0.41
2:B:163:ASN:O	2:B:166:SER:HB2	2.19	0.41
1:A:495:ASN:HB3	1:A:498:ASN:HB2	2.01	0.41
1:A:693:LEU:HD12	1:A:700:ILE:HD13	2.02	0.41
1:A:869:ILE:HG13	1:A:949:ARG:HH12	1.85	0.41
1:A:389:LYS:HG3	1:A:390:THR:HG23	2.02	0.41
1:A:847:LYS:HD2	1:A:847:LYS:N	2.36	0.41
1:A:916:GLN:OE1	2:B:66:PRO:HG3	2.21	0.41
2:B:139:PHE:HE1	2:B:247:TYR:HA	1.86	0.41
2:B:191:PRO:O	2:B:206:ASN:OD1	2.39	0.41
2:B:198:THR:OG1	2:B:199:TYR:N	2.50	0.41
1:A:59:LEU:HB2	1:A:63:GLU:HB3	2.02	0.41
1:A:148:LEU:HA	1:A:148:LEU:HD23	1.73	0.41
1:A:294:THR:OG1	1:A:296:ILE:HG22	2.20	0.41
1:A:532:THR:HA	1:A:541:PRO:HA	2.01	0.41
1:A:583:PHE:HB3	1:A:590:PHE:HD2	1.84	0.41
2:B:188:PRO:CB	2:B:210:LEU:HD11	2.51	0.41
1:A:205:ILE:CG2	1:A:266:ALA:HB3	2.51	0.41
1:A:827:ASP:O	1:A:831:SER:CB	2.66	0.41
2:B:228:ILE:HG22	2:B:264:PHE:CD1	2.56	0.41
1:A:126:ILE:HD13	1:A:126:ILE:HA	1.78	0.41
1:A:470:LYS:O	1:A:474:VAL:HG22	2.21	0.41
1:A:942:GLN:HA	1:A:945:ASP:HB3	2.03	0.41
1:A:389:LYS:CD	1:A:629:THR:HG21	2.51	0.41
1:A:394:THR:HG22	1:A:396:ASN:H	1.86	0.41
1:A:726:VAL:CG2	1:A:737:LEU:HA	2.51	0.41
1:A:900:VAL:HG13	1:A:901:ALA:N	2.36	0.41
2:B:278:LYS:HG2	2:B:295:ARG:HG2	2.03	0.41
1:A:389:LYS:HG2	1:A:629:THR:HG21	2.02	0.40
1:A:527:LEU:HD23	1:A:527:LEU:HA	1.93	0.40
1:A:1000:TYR:HA	1:A:1003:VAL:HG23	2.03	0.40
1:A:224:LYS:O	1:A:262:LEU:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:THR:O	1:A:167:THR:HG22	2.22	0.40
1:A:307:VAL:HA	1:A:310:VAL:HG22	2.04	0.40
1:A:729:ASP:HB2	1:A:749:ALA:O	2.22	0.40
2:B:21:LYS:HD2	2:B:21:LYS:HA	1.84	0.40
2:B:251:LEU:HD23	2:B:251:LEU:HA	1.92	0.40
1:A:330:LEU:H	1:A:330:LEU:HG	1.71	0.40
1:A:384:ILE:HD12	1:A:717:CYS:HB3	2.04	0.40
2:B:25:LEU:HD12	2:B:25:LEU:HA	1.82	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:484:LYS:NZ	1:A:671:GLU:OE2[3_554]	2.13	0.07

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	984/986 (100%)	920 (94%)	64 (6%)	0	100	100
2	B	286/324 (88%)	248 (87%)	38 (13%)	0	100	100
All	All	1270/1310 (97%)	1168 (92%)	102 (8%)	0	100	100

There are no Ramachandran outliers to report.

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	841/841 (100%)	830 (99%)	11 (1%)	69	82
2	B	257/286 (90%)	255 (99%)	2 (1%)	81	89
All	All	1098/1127 (97%)	1085 (99%)	13 (1%)	71	83

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	MET
1	A	58	ARG
1	A	186	ASP
1	A	328	TYR
1	A	445	PHE
1	A	515	PHE
1	A	637	LYS
1	A	641	LYS
1	A	670	ARG
1	A	961	LEU
1	A	983	TYR
2	B	270	ASP
2	B	286	TYR

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

Mol	Chain	Res	Type
1	A	71	ASN
1	A	698	GLN
1	A	795	ASN
1	A	958	GLN
1	A	999	GLN
2	B	79	GLN

5.3.3 RNA ⓘ

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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
5	NAG	B	401	2	14,14,15	0.99	1 (7%)	17,19,21	0.77	1 (5%)
3	ALF	A	1101	-	0,4,4	-	-	-		

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
5	NAG	B	401	2	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	401	NAG	O5-C1	3.47	1.49	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	401	NAG	C1-O5-C5	2.50	115.58	112.19

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	401	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1101	ALF	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	986/986 (100%)	0.03	39 (3%) 38 36	13, 34, 96, 119	0
2	B	288/324 (88%)	-0.07	7 (2%) 59 56	20, 45, 86, 114	0
All	All	1274/1310 (97%)	0.00	46 (3%) 42 40	13, 38, 95, 119	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	202	THR	8.0
2	B	164	ASP	7.1
1	A	425	GLN	6.0
1	A	404	TRP	5.8
1	A	661	ASN	5.6
1	A	494	PHE	5.5
1	A	556	GLU	4.6
1	A	552	THR	4.1
2	B	20	GLU	4.0
1	A	555	MET	3.9
2	B	23	GLU	3.7
1	A	54	LEU	3.7
1	A	408	GLN	3.7
1	A	559	GLY	3.6
1	A	554	TYR	3.4
1	A	668	ASN	3.4
2	B	165	GLU	3.3
1	A	424	ASP	3.3
1	A	600	LEU	3.3
1	A	575	GLU	3.0
1	A	551	HIS	2.9
2	B	253	PRO	2.9
1	A	100	PRO	2.9
1	A	841	SER	2.9

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Mol	Chain	Res	Type	RSRZ
2	B	21	LYS	2.9
1	A	662	ILE	2.8
1	A	167	THR	2.7
1	A	138	SER	2.6
1	A	98	GLN	2.5
1	A	52	MET	2.5
1	A	181	ALA	2.5
1	A	652	THR	2.5
1	A	496	SER	2.5
1	A	292	GLU	2.4
1	A	426	SER	2.4
1	A	557	LEU	2.4
1	A	560	LEU	2.3
1	A	407	ASN	2.3
1	A	599	GLY	2.3
1	A	99	THR	2.3
1	A	601	LEU	2.2
1	A	405	PHE	2.2
1	A	403	LEU	2.2
1	A	410	PHE	2.2
1	A	51	GLY	2.1
1	A	581	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

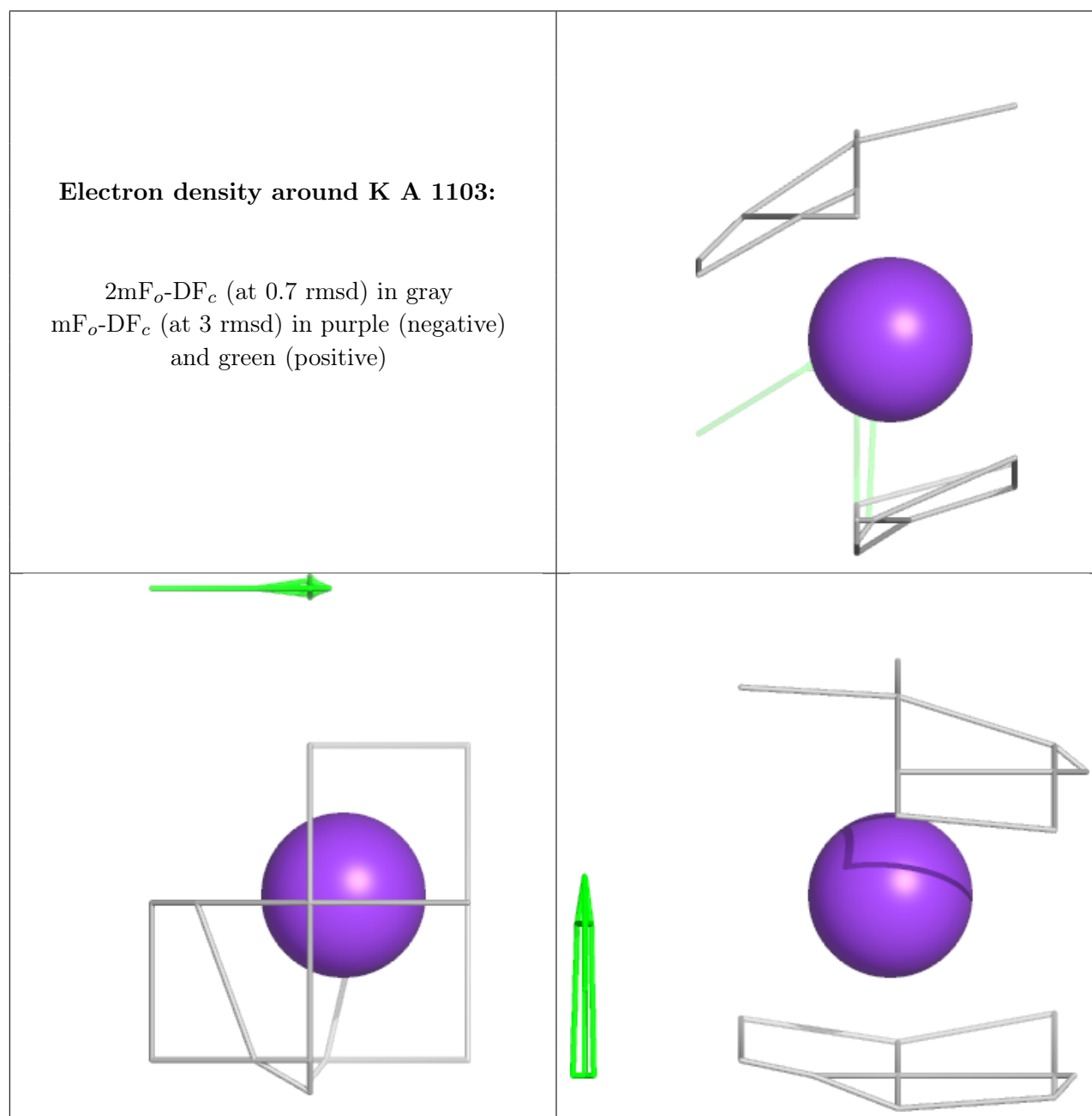
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	NAG	B	401	14/15	0.66	0.44	74,85,93,94	0

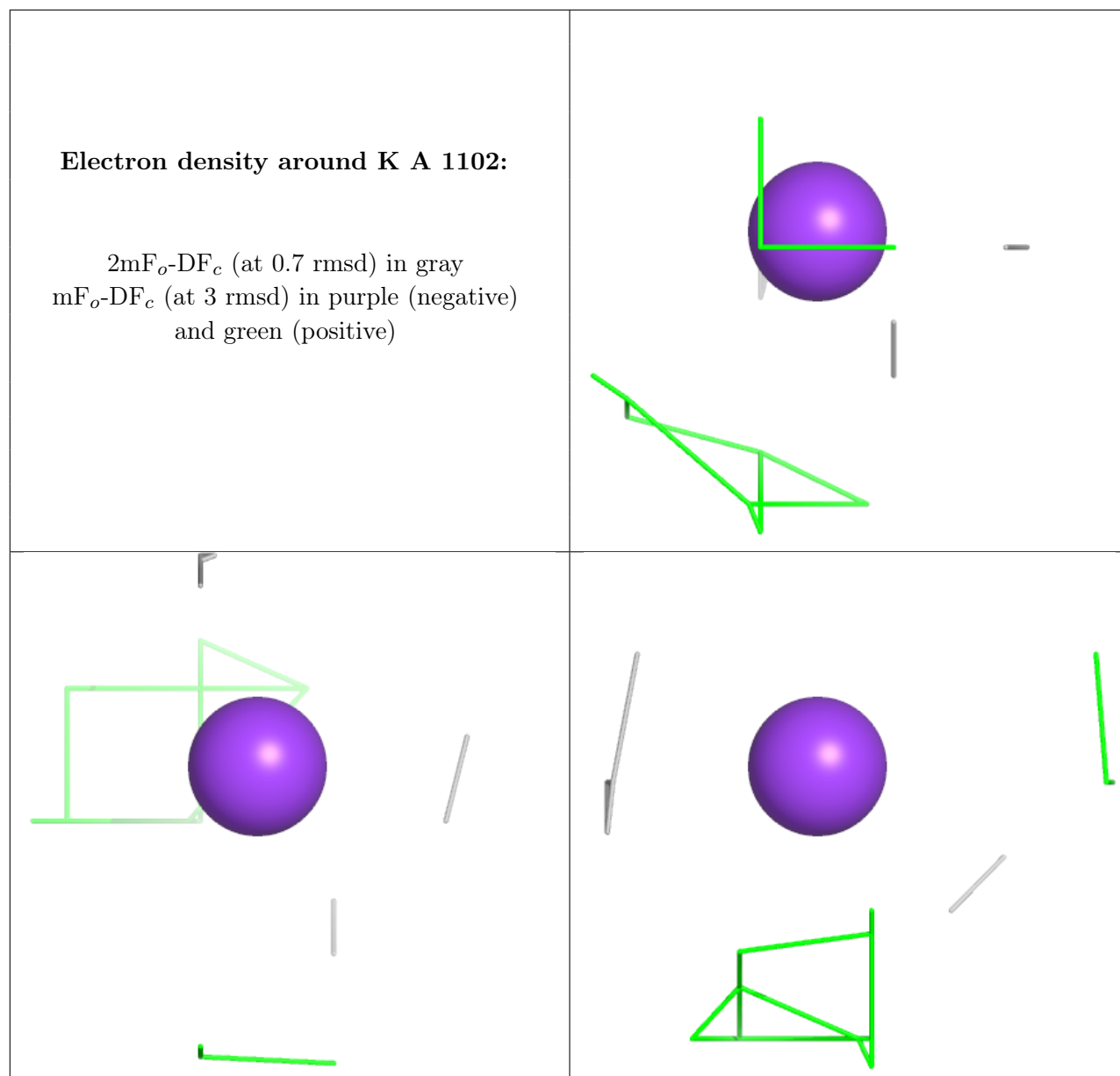
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	K	A	1103	1/1	0.84	0.15	24,24,24,24	0
4	K	A	1102	1/1	0.88	0.15	23,23,23,23	0
3	ALF	A	1101	5/5	0.93	0.27	19,21,25,26	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





6.5 Other polymers ⓘ

There are no such residues in this entry.