



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 28, 2022 – 04:16 PM EDT

PDB ID : 7N93  
Title : P70 S6K1 IN COMPLEX WITH MSC2363318A-1  
Authors : Mochalkin, I.  
Deposited on : 2021-06-16  
Resolution : 2.74 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at <http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.29  
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) : 2.29

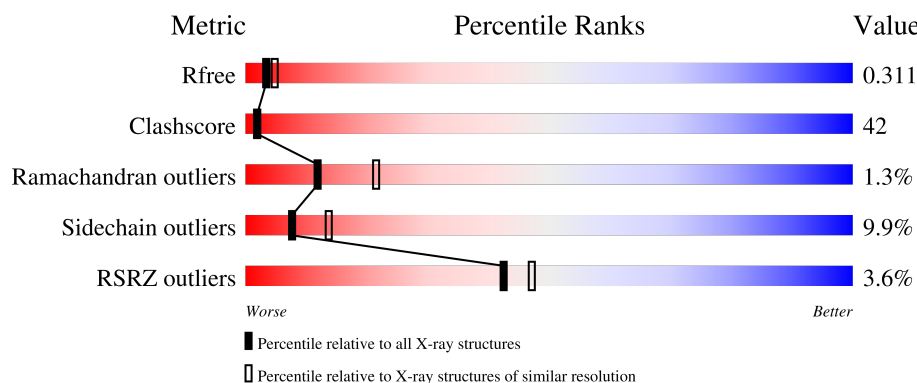
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.74 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	1271 (2.76-2.72)
Clashscore	141614	1322 (2.76-2.72)
Ramachandran outliers	138981	1297 (2.76-2.72)
Sidechain outliers	138945	1298 (2.76-2.72)
RSRZ outliers	127900	1243 (2.76-2.72)

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

Mol	Chain	Length	Quality of chain
1	A	340	<div> <div>5%</div> <div>45%</div> <div>39%</div> <div>8%</div> <div>8%</div> </div>
1	B	340	<div> <div>5%</div> <div>43%</div> <div>41%</div> <div>6%</div> <div>9%</div> </div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5062 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.

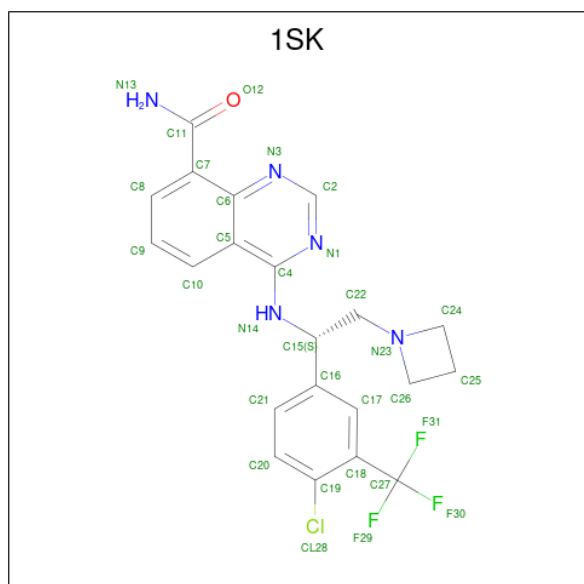
- Molecule 1 is a protein called Ribosomal protein S6 kinase beta-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	312	Total	C	N	O	P	S	0	2	0
			2515	1613	432	453	2	15			
1	B	309	Total	C	N	O	P	S	0	1	0
			2481	1597	427	442	1	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	412	GLU	THR	engineered mutation	UNP P23443
B	412	GLU	THR	engineered mutation	UNP P23443

- Molecule 2 is 4-((1S)-2-(azetidin-1-yl)-1-[4-chloro-3-(trifluoromethyl)phenyl]ethyl)amino)quinazoline-8-carboxamide (three-letter code: 1SK) (formula: C<sub>21</sub>H<sub>19</sub>ClF<sub>3</sub>N<sub>5</sub>O) (labeled as "Ligand of Interest" by depositor).



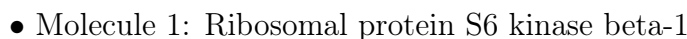
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 31	C 21	Cl 1	F 3	N 5	O 1	0	0
2	B	1	Total 31	C 21	Cl 1	F 3	N 5	O 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total 3	O 3	0	0
3	B	1	Total 1	O 1	0	0



- Molecule 1: Ribosomal protein S6 kinase beta-1



SEP	PRO	ASP	ASP	SER	THR	LEU	SER	GLU	SER	ALA	N405	Q406	V407	F408	L409	G410	F411	E412	Y413	V414	A415	P416	S417	V418	LEU	GLU	SER
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.02Å 75.45Å 151.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.74 47.53 – 2.74	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.74) 98.9 (47.53-2.74)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.27 (at 2.73Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, $R_{free}$	0.211 , 0.290 0.242 , 0.311	Depositor DCC
$R_{free}$ test set	944 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	69.8	Xtriage
Anisotropy	0.470	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 26.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	5062	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TPO, SEP, 1SK

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.44	0/2547	0.65	3/3423 (0.1%)
1	B	0.45	0/2523	0.76	11/3393 (0.3%)
All	All	0.45	0/5070	0.71	14/6816 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	384	SER	CB-CA-C	-8.96	93.07	110.10
1	B	298	ARG	CB-CA-C	-8.96	92.48	110.40
1	B	131	VAL	CB-CA-C	-8.85	94.58	111.40
1	B	298	ARG	N-CA-C	8.31	133.44	111.00
1	B	377	GLU	CB-CA-C	-8.05	94.30	110.40
1	A	296	GLU	CB-CA-C	7.86	126.12	110.40
1	B	136	ASP	CB-CA-C	-7.67	95.06	110.40
1	A	131	VAL	CB-CA-C	-7.41	97.32	111.40
1	A	384	SER	CB-CA-C	-7.18	96.47	110.10
1	B	406	GLN	N-CA-C	-6.89	92.39	111.00
1	B	320	ALA	CB-CA-C	-6.69	100.07	110.10
1	B	136	ASP	N-CA-C	6.53	128.63	111.00
1	B	331	ASN	CB-CA-C	6.36	123.11	110.40
1	B	129	MET	CB-CA-C	-5.46	99.47	110.40



There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	384	SER	Peptide

## 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	2515	0	2533	220	0
1	B	2481	0	2515	203	0
2	A	31	0	0	3	0
2	B	31	0	0	2	0
3	A	3	0	0	0	0
3	B	1	0	0	0	0
All	All	5062	0	5048	422	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 42.

All (422) 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:330:ARG:HH11	1:A:330:ARG:CB	1.03	1.61
1:A:330:ARG:CB	1:A:330:ARG:NH1	1.83	1.41
1:A:417:SER:O	1:A:418:VAL:HG22	1.22	1.37
1:B:304:LYS:HD2	1:B:308:CYS:SG	1.69	1.33
1:A:330:ARG:NH1	1:A:330:ARG:HB2	1.41	1.26
1:A:406:GLN:HE21	1:A:408:PHE:N	1.33	1.23
1:A:241:LYS:C	1:A:242:GLU:OE1	1.84	1.15
1:A:242:GLU:OE1	1:A:242:GLU:N	1.79	1.15
1:B:319:GLU:OE1	1:B:319:GLU:N	1.80	1.14
1:B:308:CYS:O	1:B:330:ARG:NH2	1.83	1.11
1:B:409:LEU:HD12	1:B:410:GLY:H	1.12	1.08
1:A:417:SER:O	1:A:418:VAL:CG2	2.01	1.08
1:A:330:ARG:NH1	1:A:330:ARG:HB3	1.64	1.07
1:A:99:LYS:HE2	1:A:104:LYS:HE2	1.35	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:VAL:C	1:B:116:THR:HG23	1.75	1.05
1:A:229[A]:GLN:NE2	1:B:358:GLU:HG3	1.71	1.05
1:A:321:ARG:CB	1:A:321:ARG:HH11	1.69	1.04
1:B:127:LYS:HG3	1:B:166:GLY:O	1.55	1.03
1:A:87:ARG:HB3	1:A:88:PRO:HD2	1.36	1.03
1:A:132:ARG:HB2	1:A:132:ARG:HH11	1.22	1.02
1:B:131:VAL:HA	1:B:137:THR:HG21	1.39	1.02
1:A:321:ARG:HH11	1:A:321:ARG:CG	1.71	1.01
1:B:409:LEU:HD12	1:B:410:GLY:N	1.76	1.00
1:B:304:LYS:CD	1:B:308:CYS:SG	2.52	0.98
1:A:191:MET:HE2	1:A:191:MET:HA	1.46	0.98
1:B:131:VAL:HA	1:B:137:THR:CG2	1.94	0.97
1:A:406:GLN:HE21	1:A:408:PHE:H	0.98	0.97
1:A:199:LEU:HD23	1:A:279:LEU:HD11	1.45	0.95
1:A:132:ARG:HB2	1:A:132:ARG:NH1	1.80	0.95
1:A:406:GLN:NE2	1:A:408:PHE:N	2.15	0.95
1:A:167:LYS:NZ	1:A:394:SEP:O	1.99	0.95
1:A:260:MET:CE	1:A:264:ILE:HG21	1.96	0.94
1:B:297:ASN:ND2	1:B:300:LYS:HB2	1.83	0.93
1:B:349:HIS:ND1	1:B:350:PRO:HD2	1.85	0.91
1:B:114:ALA:O	1:B:115:ASN:HB2	1.71	0.90
1:A:313:PRO:HD2	1:A:316:LEU:HD12	1.51	0.90
1:A:113:GLY:O	1:A:116:THR:OG1	1.91	0.89
1:B:159:ILE:HG22	1:B:160:TYR:CD1	2.07	0.89
1:B:405:ASN:C	1:B:405:ASN:HD22	1.74	0.89
1:B:303:ASP:O	1:B:307:LYS:HG3	1.73	0.89
1:A:114:ALA:O	1:A:115:ASN:HB2	1.70	0.88
1:A:163:GLN:O	1:A:164:THR:HG23	1.74	0.88
1:A:406:GLN:NE2	1:A:408:PHE:H	1.70	0.88
1:B:405:ASN:ND2	1:B:405:ASN:O	2.07	0.87
1:A:141:LYS:NZ	1:A:407:VAL:O	2.07	0.86
1:B:306:LEU:O	1:B:330:ARG:NH1	2.09	0.86
1:A:132:ARG:NH1	1:A:132:ARG:CB	2.37	0.86
1:B:103:GLY:HA2	1:B:126:LYS:HZ1	1.41	0.85
1:B:406:GLN:HA	1:B:406:GLN:OE1	1.74	0.85
1:B:297:ASN:HD21	1:B:300:LYS:HB2	1.42	0.84
1:A:87:ARG:HB3	1:A:88:PRO:CD	2.06	0.84
1:A:297:ASN:C	1:A:297:ASN:HD22	1.79	0.84
1:B:297:ASN:HD21	1:B:300:LYS:CB	1.89	0.84
1:A:321:ARG:CB	1:A:321:ARG:NH1	2.41	0.84
1:B:160:TYR:CE2	1:B:414:VAL:HG13	2.13	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:358:GLU:HA	1:B:358:GLU:OE1	1.78	0.83
1:A:321:ARG:NH1	1:A:321:ARG:HB3	1.92	0.82
1:A:330:ARG:HH11	1:A:330:ARG:HB2	0.65	0.82
1:A:366:GLU:OE1	1:A:366:GLU:HA	1.79	0.82
1:A:119:ILE:HG13	1:A:376:GLU:HG2	1.62	0.81
1:A:191:MET:HA	1:A:191:MET:CE	2.09	0.81
1:A:321:ARG:HH11	1:A:321:ARG:HG3	1.44	0.81
1:B:296:GLU:N	1:B:296:GLU:OE1	2.13	0.81
1:A:127:LYS:HG3	1:A:166:GLY:O	1.81	0.80
1:A:297:ASN:HD21	1:A:300:LYS:H	1.28	0.80
1:B:111:VAL:HA	1:B:116:THR:HG23	1.64	0.80
1:B:130:ILE:HG22	1:B:137:THR:HG22	1.64	0.79
1:A:127:LYS:HG3	1:A:166:GLY:C	2.03	0.79
1:A:150:VAL:HG12	1:A:150:VAL:O	1.80	0.79
1:A:342:ASP:OD2	1:A:343:ALA:N	2.16	0.78
1:A:102:TYR:O	1:A:126:LYS:HE3	1.84	0.78
1:A:297:ASN:ND2	1:A:300:LYS:H	1.82	0.77
1:B:304:LYS:HD2	1:B:308:CYS:HG	1.47	0.76
1:B:111:VAL:O	1:B:116:THR:HG23	1.84	0.76
1:B:133:ASN:C	1:B:133:ASN:OD1	2.24	0.76
1:B:186:ARG:O	1:B:187:GLU:OE2	2.03	0.76
1:A:218:ASP:OD2	1:A:220:LYS:HE2	1.86	0.76
1:B:381:GLN:HE21	1:B:381:GLN:N	1.85	0.75
1:A:189:ILE:CD1	1:A:289:GLY:N	2.49	0.75
1:A:98:GLY:HA2	1:A:386:PHE:CE1	2.21	0.75
1:B:183:GLN:OE1	1:B:183:GLN:HA	1.87	0.75
1:A:260:MET:HE1	1:A:264:ILE:HG21	1.68	0.75
1:B:99:LYS:HG3	1:B:104:LYS:HG2	1.69	0.74
1:A:87:ARG:CB	1:A:88:PRO:HD2	2.16	0.74
1:B:159:ILE:HG22	1:B:160:TYR:CE1	2.23	0.74
1:A:373:LEU:HD23	1:A:378:ASP:HB3	1.69	0.73
1:A:394:SEP:O	1:A:394:SEP:O3P	2.05	0.73
1:B:297:ASN:HD21	1:B:300:LYS:CG	2.02	0.73
1:A:199:LEU:HD23	1:A:279:LEU:CD1	2.17	0.73
1:B:381:GLN:NE2	1:B:381:GLN:HA	2.04	0.72
1:A:132:ARG:HH11	1:A:132:ARG:CB	2.00	0.72
1:A:296:GLU:HB3	1:A:300:LYS:HD2	1.70	0.72
1:B:349:HIS:CG	1:B:350:PRO:HD2	2.25	0.72
1:A:94:LEU:C	1:A:95:ARG:HG2	2.10	0.72
1:A:417:SER:C	1:A:418:VAL:HG22	2.10	0.72
1:A:94:LEU:O	1:A:95:ARG:HG2	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:297:ASN:C	1:B:297:ASN:HD22	1.93	0.71
1:A:176:SER:HB2	1:A:373:LEU:HD21	1.73	0.70
1:A:296:GLU:HB3	1:A:300:LYS:CD	2.21	0.70
1:B:405:ASN:C	1:B:405:ASN:ND2	2.45	0.70
1:B:192:GLU:OE1	1:B:317:THR:HG23	1.91	0.70
1:B:98:GLY:HA2	1:B:386:PHE:CG	2.27	0.70
1:B:297:ASN:HD21	1:B:300:LYS:HG3	1.57	0.70
1:B:194:THR:HG23	1:B:368:PRO:HG3	1.73	0.69
1:A:257:ILE:HD13	1:A:257:ILE:H	1.56	0.69
1:B:159:ILE:CG2	1:B:160:TYR:CE1	2.75	0.69
1:B:303:ASP:HB3	1:B:307:LYS:HD3	1.75	0.69
1:A:241:LYS:O	1:A:241:LYS:HG2	1.93	0.69
1:A:321:ARG:CG	1:A:321:ARG:NH1	2.44	0.68
1:A:132:ARG:HB3	1:A:132:ARG:CZ	2.22	0.68
1:A:297:ASN:C	1:A:297:ASN:ND2	2.47	0.68
1:A:406:GLN:HG2	1:A:407:VAL:N	2.09	0.68
1:B:103:GLY:HA2	1:B:126:LYS:NZ	2.07	0.68
1:B:111:VAL:O	1:B:116:THR:CG2	2.41	0.68
1:B:225:MET:HG3	1:B:235:THR:HG21	1.76	0.68
1:B:111:VAL:CA	1:B:116:THR:HG23	2.24	0.67
1:B:303:ASP:O	1:B:307:LYS:CG	2.40	0.67
1:B:94:LEU:O	1:B:95:ARG:HG2	1.94	0.67
1:B:303:ASP:C	1:B:307:LYS:HD3	2.15	0.67
1:B:381:GLN:NE2	1:B:381:GLN:CA	2.58	0.67
1:A:407:VAL:O	1:A:407:VAL:HG22	1.93	0.67
1:A:418:VAL:HG23	1:A:418:VAL:O	1.95	0.67
1:A:103:GLY:HA2	1:A:126:LYS:HE2	1.77	0.66
1:B:149:GLU:C	1:B:149:GLU:OE1	2.33	0.66
1:B:194:THR:HG23	1:B:368:PRO:CG	2.25	0.66
1:B:130:ILE:CG2	1:B:137:THR:HA	2.24	0.66
1:B:380:SER:C	1:B:381:GLN:HE21	1.97	0.66
1:A:164:THR:O	1:A:165:GLY:C	2.33	0.66
1:B:349:HIS:CE1	1:B:350:PRO:HD2	2.29	0.66
1:B:87:ARG:HB3	1:B:88:PRO:HD2	1.76	0.66
1:B:99:LYS:CG	1:B:104:LYS:HG2	2.25	0.66
1:B:98:GLY:HA2	1:B:386:PHE:CD2	2.32	0.65
1:A:191:MET:CE	1:A:191:MET:CA	2.74	0.65
1:B:222:GLU:O	2:B:501:ISK:C26	2.44	0.65
1:B:93:LEU:CD2	1:B:122:MET:HE1	2.27	0.64
1:A:417:SER:O	1:A:418:VAL:CB	2.44	0.64
1:B:303:ASP:HB3	1:B:307:LYS:CD	2.27	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:GLY:HA2	1:B:386:PHE:CD1	2.32	0.64
1:B:111:VAL:HA	1:B:116:THR:CG2	2.27	0.64
1:B:271:ASN:O	1:B:274:VAL:HG13	1.97	0.64
1:B:297:ASN:ND2	1:B:297:ASN:C	2.50	0.64
1:B:381:GLN:HE21	1:B:381:GLN:CA	2.10	0.64
1:B:160:TYR:CD2	1:B:414:VAL:HG13	2.32	0.64
1:B:407:VAL:O	1:B:407:VAL:HG22	1.97	0.64
1:A:132:ARG:CB	1:A:132:ARG:CZ	2.75	0.63
1:A:169:TYR:O	1:A:170:LEU:HD23	1.98	0.63
1:B:349:HIS:ND1	1:B:350:PRO:CD	2.61	0.63
1:A:384:SER:O	1:A:388:ARG:HB2	1.99	0.63
1:A:303:ASP:O	1:A:307:LYS:HG2	1.99	0.63
1:A:131:VAL:HG23	1:A:132:ARG:N	2.12	0.62
1:A:189:ILE:HD11	1:A:289:GLY:N	2.14	0.62
1:A:260:MET:HE3	1:A:264:ILE:HG21	1.80	0.62
1:B:94:LEU:C	1:B:95:ARG:HG2	2.19	0.62
1:A:229[A]:GLN:NE2	1:B:358:GLU:CG	2.56	0.62
1:A:133:ASN:ND2	1:A:136:ASP:H	1.97	0.62
1:B:218:ASP:OD2	1:B:260:MET:HE3	2.00	0.62
1:A:97:LEU:HD11	1:A:107:GLN:HB2	1.82	0.61
1:B:130:ILE:HG22	1:B:137:THR:HA	1.81	0.61
1:A:201:GLU:OE1	1:A:232:VAL:HG13	1.99	0.61
1:A:321:ARG:NH1	1:A:321:ARG:HG3	2.11	0.61
1:B:265:LEU:HD13	1:B:302:ILE:HG12	1.81	0.61
1:A:133:ASN:HD22	1:A:136:ASP:CG	2.04	0.61
1:A:347:GLN:HB3	1:A:357:TRP:CZ2	2.34	0.61
1:A:373:LEU:CD2	1:A:378:ASP:HB3	2.30	0.61
1:A:144:ARG:HD2	1:A:144:ARG:O	2.01	0.60
1:B:373:LEU:HD22	1:B:378:ASP:HB3	1.83	0.60
1:A:417:SER:O	1:A:418:VAL:HG13	2.01	0.60
1:A:102:TYR:CE1	1:A:129:MET:HB2	2.37	0.60
1:A:366:GLU:OE1	1:A:367:PRO:HD2	2.02	0.60
1:B:362:ALA:O	1:B:363:ARG:HB2	2.01	0.60
1:A:132:ARG:NH1	1:A:132:ARG:HB3	2.17	0.60
1:B:99:LYS:HG3	1:B:103:GLY:O	2.01	0.60
1:A:271:ASN:OD1	1:A:273:ALA:N	2.26	0.59
1:B:159:ILE:CG2	1:B:160:TYR:CD1	2.84	0.59
1:A:102:TYR:CD1	1:A:129:MET:HB2	2.37	0.59
1:A:313:PRO:HB2	1:A:315:TYR:CE2	2.36	0.59
1:B:87:ARG:HB3	1:B:88:PRO:CD	2.33	0.59
1:B:103:GLY:CA	1:B:126:LYS:HZ1	2.14	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:297:ASN:ND2	1:B:300:LYS:HG3	2.17	0.59
1:A:260:MET:CE	1:A:264:ILE:CG2	2.78	0.59
1:A:317:THR:O	1:A:321:ARG:HG2	2.02	0.59
1:A:313:PRO:CB	1:A:315:TYR:CE2	2.86	0.58
1:A:222:GLU:O	2:A:501:ISK:C25	2.51	0.58
1:B:94:LEU:HB2	1:B:107:GLN:HG2	1.84	0.58
1:A:91:PHE:HB3	1:A:108:VAL:HG21	1.85	0.58
1:A:163:GLN:O	1:A:164:THR:CG2	2.49	0.58
1:B:415:ALA:O	1:B:418:VAL:HG23	2.03	0.58
1:A:242:GLU:OE1	1:A:242:GLU:CA	2.52	0.58
1:B:222:GLU:O	2:B:501:ISK:C25	2.51	0.58
1:B:161:ALA:O	1:B:412:GLU:HA	2.04	0.58
1:B:377:GLU:O	1:B:377:GLU:HG2	1.99	0.58
1:A:343:ALA:O	1:A:347:GLN:HG3	2.03	0.58
1:A:384:SER:HA	1:A:387:THR:OG1	2.04	0.58
1:B:127:LYS:HG3	1:B:166:GLY:C	2.23	0.58
1:B:92:GLU:HG3	1:B:92:GLU:O	2.04	0.57
1:A:163:GLN:NE2	1:A:412:GLU:OE2	2.37	0.57
1:B:87:ARG:CB	1:B:88:PRO:CD	2.83	0.57
1:B:266:MET:O	1:B:267:ARG:C	2.40	0.57
1:A:189:ILE:HD11	1:A:288:THR:C	2.25	0.57
1:A:103:GLY:HA3	1:A:124:VAL:O	2.05	0.57
1:B:277:TRP:O	1:B:280:GLY:N	2.38	0.57
1:A:198:TYR:O	1:A:202:ILE:HG13	2.05	0.57
1:A:406:GLN:HG2	1:A:407:VAL:H	1.69	0.57
1:B:282:LEU:O	1:B:286:MET:HG3	2.05	0.56
1:B:326:LYS:HB3	1:B:336:LEU:HB2	1.87	0.56
1:B:149:GLU:OE1	1:B:149:GLU:O	2.23	0.56
1:B:358:GLU:OE1	1:B:358:GLU:CA	2.52	0.56
1:A:98:GLY:HA2	1:A:386:PHE:CD1	2.41	0.56
1:B:285:ASP:HA	1:B:290:ALA:O	2.05	0.56
1:A:294:THR:HG22	1:A:295:GLY:HA3	1.88	0.56
1:A:176:SER:O	1:A:381:GLN:HG3	2.06	0.56
1:A:130:ILE:HD13	1:A:130:ILE:N	2.20	0.55
1:A:409:LEU:O	1:A:410:GLY:C	2.43	0.55
1:B:163:GLN:HG3	1:B:163:GLN:O	2.05	0.55
1:B:192:GLU:OE1	1:B:317:THR:CG2	2.55	0.55
1:B:330:ARG:H	1:B:330:ARG:HD2	1.72	0.55
1:A:356:ASN:OD1	1:A:356:ASN:O	2.24	0.55
1:A:102:TYR:CD1	1:A:130:ILE:CD1	2.90	0.54
1:A:241:LYS:CA	1:A:242:GLU:OE1	2.55	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:ILE:HG23	1:A:298:ARG:HH21	1.72	0.54
1:B:339:GLY:HA3	1:B:345:GLU:OE2	2.07	0.54
1:B:93:LEU:HD21	1:B:122:MET:CE	2.38	0.54
1:A:417:SER:O	1:A:418:VAL:CG1	2.55	0.54
1:B:384:SER:HA	1:B:387:THR:OG1	2.08	0.54
1:B:93:LEU:CD2	1:B:122:MET:CE	2.85	0.53
1:B:229:GLN:O	1:B:367:PRO:HG2	2.07	0.53
1:B:98:GLY:HA2	1:B:386:PHE:CE1	2.44	0.52
1:A:113:GLY:O	1:A:116:THR:CB	2.57	0.52
1:A:152:HIS:CG	1:A:153:PRO:HD2	2.44	0.52
1:B:98:GLY:HA2	1:B:386:PHE:CE2	2.44	0.52
1:A:156:VAL:HB	1:A:235:THR:HG22	1.91	0.52
1:A:339:GLY:HA3	1:A:345:GLU:OE2	2.09	0.52
2:A:501:1SK:N3	2:A:501:1SK:N13	2.56	0.52
1:A:199:LEU:CD2	1:A:279:LEU:CD1	2.88	0.52
1:B:130:ILE:HG22	1:B:137:THR:CG2	2.38	0.52
1:B:323:LEU:HG	1:B:327:LEU:CD1	2.39	0.52
1:A:313:PRO:HD2	1:A:316:LEU:CD1	2.32	0.52
1:B:191:MET:O	1:B:192:GLU:C	2.49	0.51
1:B:303:ASP:O	1:B:307:LYS:HD3	2.11	0.51
1:A:189:ILE:HD11	1:A:288:THR:CA	2.40	0.51
1:A:406:GLN:HE21	1:A:407:VAL:C	2.08	0.51
1:B:384:SER:O	1:B:388:ARG:HG3	2.10	0.51
1:B:262:PRO:HG3	1:B:305:ILE:CG2	2.41	0.51
1:B:377:GLU:O	1:B:377:GLU:CG	2.52	0.51
1:B:218:ASP:HB2	1:B:260:MET:HE2	1.91	0.51
1:B:103:GLY:CA	1:B:126:LYS:NZ	2.73	0.51
1:B:111:VAL:O	1:B:111:VAL:HG12	2.11	0.51
1:A:377:GLU:OE2	1:B:212:LYS:NZ	2.44	0.51
1:B:201:GLU:OE1	1:B:232:VAL:HG13	2.11	0.51
1:B:94:LEU:O	1:B:95:ARG:CG	2.59	0.51
1:A:358[A]:GLU:HA	1:A:358[A]:GLU:OE2	2.11	0.50
1:B:297:ASN:ND2	1:B:300:LYS:CB	2.56	0.50
1:B:330:ARG:HD2	1:B:330:ARG:N	2.26	0.50
1:A:148:GLU:HG3	1:A:413:TYR:CE1	2.47	0.50
1:B:93:LEU:HD21	1:B:122:MET:HE1	1.92	0.50
1:B:103:GLY:HA2	1:B:126:LYS:CE	2.42	0.50
1:A:366:GLU:OE1	1:A:367:PRO:CD	2.59	0.50
1:A:127:LYS:HD2	1:A:166:GLY:C	2.32	0.49
1:A:114:ALA:O	1:A:115:ASN:CB	2.49	0.49
1:B:228:HIS:H	1:B:228:HIS:CD2	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304:LYS:HD2	1:A:308:CYS:SG	2.52	0.49
1:B:159:ILE:HG21	1:B:160:TYR:CE1	2.45	0.49
1:A:406:GLN:HG2	1:A:408:PHE:H	1.77	0.49
1:A:229[A]:GLN:O	1:A:367:PRO:HG2	2.13	0.49
1:B:150:VAL:HG11	1:B:155:ILE:HD13	1.95	0.49
1:A:102:TYR:O	1:A:126:LYS:CE	2.58	0.49
1:A:191:MET:HE2	1:A:191:MET:CA	2.28	0.49
1:B:161:ALA:HB3	1:B:413:TYR:HB3	1.94	0.49
1:A:204:MET:HE2	1:A:361:LEU:HD13	1.93	0.49
1:A:356:ASN:OD1	1:A:356:ASN:C	2.51	0.49
1:A:189:ILE:CD1	1:A:288:THR:C	2.80	0.49
1:A:204:MET:CE	1:A:361:LEU:HD13	2.42	0.49
1:A:114:ALA:C	1:A:116:THR:H	2.15	0.49
1:A:257:ILE:H	1:A:257:ILE:CD1	2.26	0.49
1:A:358[A]:GLU:OE2	1:A:358[A]:GLU:CA	2.61	0.49
1:B:271:ASN:O	1:B:273:ALA:N	2.45	0.49
1:B:353:ARG:HG3	1:B:354:HIS:CD2	2.47	0.49
1:A:124:VAL:HG22	1:A:169:TYR:CD1	2.48	0.48
1:B:114:ALA:O	1:B:115:ASN:CB	2.50	0.48
1:B:293:PHE:O	1:B:301:THR:HG23	2.12	0.48
1:B:99:LYS:HD2	1:B:104:LYS:HE2	1.95	0.48
1:B:160:TYR:HE2	1:B:414:VAL:HG13	1.75	0.48
1:A:131:VAL:CG2	1:A:132:ARG:N	2.77	0.48
1:A:316:LEU:HB3	1:A:321:ARG:CD	2.43	0.48
1:B:169:TYR:O	1:B:170:LEU:HD23	2.13	0.48
1:A:191:MET:O	1:A:194:THR:HB	2.14	0.48
1:B:303:ASP:CB	1:B:307:LYS:CD	2.90	0.48
1:A:216:TYR:O	1:A:217:ARG:HB2	2.14	0.48
1:A:295:GLY:O	1:A:296:GLU:C	2.52	0.48
1:B:303:ASP:O	1:B:307:LYS:CD	2.61	0.48
1:A:99:LYS:HG2	1:A:104:LYS:HG2	1.95	0.48
1:B:93:LEU:HD23	1:B:122:MET:HE1	1.95	0.48
1:B:103:GLY:N	1:B:126:LYS:NZ	2.62	0.48
1:A:96:VAL:HG22	1:A:106:PHE:HE1	1.79	0.48
1:A:199:LEU:CD2	1:A:279:LEU:HG	2.44	0.48
1:B:263:GLU:OE1	1:B:274:VAL:HG12	2.14	0.48
1:B:262:PRO:HG3	1:B:305:ILE:HG22	1.95	0.48
1:A:205:ALA:O	1:A:209:LEU:HG	2.13	0.47
1:A:370:LYS:O	1:A:371:PRO:C	2.51	0.47
1:A:407:VAL:O	1:A:407:VAL:CG2	2.61	0.47
1:A:316:LEU:CB	1:A:321:ARG:HD3	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:VAL:N	1:B:137:THR:HG22	2.28	0.47
1:A:316:LEU:O	1:A:321:ARG:HD3	2.14	0.47
1:A:406:GLN:CG	1:A:407:VAL:N	2.77	0.47
1:B:130:ILE:HG22	1:B:137:THR:CA	2.45	0.47
1:B:204:MET:HE2	1:B:361:LEU:HD13	1.95	0.47
1:A:103:GLY:CA	1:A:126:LYS:HE2	2.45	0.47
1:A:126:LYS:O	1:A:130:ILE:HG12	2.14	0.47
1:A:127:LYS:HD2	1:A:167:LYS:N	2.30	0.47
1:B:296:GLU:H	1:B:296:GLU:CD	2.14	0.46
1:A:162:PHE:CE2	1:A:169:TYR:CG	3.03	0.46
1:B:94:LEU:O	1:B:95:ARG:HD3	2.14	0.46
1:B:98:GLY:HA2	1:B:386:PHE:CZ	2.50	0.46
1:A:97:LEU:O	1:A:386:PHE:HD1	1.99	0.46
1:A:207:GLY:O	1:A:211:GLN:HG3	2.15	0.46
1:B:146:ILE:HG23	1:B:214:ILE:HD13	1.98	0.46
1:B:303:ASP:CB	1:B:307:LYS:HD2	2.45	0.46
1:A:159:ILE:HG22	1:A:160:TYR:CE1	2.50	0.46
1:A:192:GLU:OE1	1:A:317:THR:HG23	2.16	0.46
1:A:102:TYR:CE1	1:A:129:MET:CB	2.99	0.46
1:B:206:LEU:HD11	1:B:219:LEU:HD11	1.97	0.46
1:B:99:LYS:CB	1:B:104:LYS:HG2	2.46	0.46
1:B:184:LEU:HA	1:B:184:LEU:HD23	1.71	0.46
1:A:102:TYR:CD1	1:A:130:ILE:HD13	2.52	0.45
1:B:99:LYS:HB2	1:B:104:LYS:HG2	1.98	0.45
1:B:103:GLY:HA2	1:B:126:LYS:HE3	1.97	0.45
1:B:144:ARG:NH2	1:B:413:TYR:HB2	2.31	0.45
1:A:282:LEU:O	1:A:282:LEU:HD22	2.17	0.45
1:B:152:HIS:HB3	1:B:155:ILE:HD12	1.97	0.45
1:A:107:GLN:NE2	1:A:379:VAL:CG1	2.80	0.45
1:B:133:ASN:OD1	1:B:134:ALA:N	2.48	0.45
1:B:391:PRO:O	1:B:392:VAL:HG23	2.17	0.45
1:A:162:PHE:CE2	1:A:169:TYR:CB	3.00	0.45
1:A:394:SEP:HA	1:A:395:PRO:HD3	1.81	0.45
1:A:97:LEU:CD1	1:A:107:GLN:HB2	2.46	0.45
1:A:307:LYS:HB2	1:A:307:LYS:HE3	1.58	0.45
1:B:163:GLN:O	1:B:163:GLN:CG	2.65	0.45
1:B:183:GLN:O	1:B:187:GLU:HG2	2.15	0.45
1:A:127:LYS:HZ1	1:A:163:GLN:HB2	1.81	0.45
1:A:170:LEU:HD21	1:A:411:PHE:CZ	2.52	0.45
1:B:271:ASN:C	1:B:273:ALA:N	2.71	0.45
1:A:182:MET:HB3	1:A:186:ARG:NH1	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:GLN:OE1	1:A:186:ARG:NH2	2.50	0.45
1:B:266:MET:O	1:B:268:SER:N	2.50	0.44
1:B:283:MET:O	1:B:284:TYR:C	2.56	0.44
1:A:127:LYS:HD2	1:A:166:GLY:CA	2.47	0.44
1:A:320:ALA:O	1:A:321:ARG:C	2.53	0.44
1:B:307:LYS:HG3	1:B:307:LYS:H	1.53	0.44
1:A:189:ILE:HD11	1:A:288:THR:N	2.32	0.44
1:A:316:LEU:CB	1:A:321:ARG:CD	2.96	0.44
1:A:228:HIS:CE1	1:A:229[A]:GLN:HG2	2.52	0.44
1:B:188:GLY:C	1:B:189:ILE:HG13	2.38	0.44
1:A:382:PHE:N	1:A:382:PHE:CD1	2.85	0.44
1:A:313:PRO:HB3	1:A:315:TYR:CE2	2.53	0.44
1:A:350:PRO:O	1:A:353:ARG:HD3	2.17	0.44
1:B:175:LEU:HD22	1:B:233:LYS:HG3	2.00	0.44
1:B:297:ASN:ND2	1:B:300:LYS:CG	2.76	0.44
1:B:406:GLN:OE1	1:B:406:GLN:CA	2.58	0.44
1:A:295:GLY:C	1:A:297:ASN:N	2.70	0.44
1:B:416:PRO:O	1:B:418:VAL:N	2.51	0.43
1:B:304:LYS:CD	1:B:308:CYS:HG	2.19	0.43
1:A:316:LEU:HB3	1:A:321:ARG:HD2	2.00	0.43
1:A:124:VAL:HA	1:A:168:LEU:O	2.18	0.43
1:A:144:ARG:NH2	1:A:413:TYR:HB2	2.34	0.43
1:A:219:LEU:HD23	1:A:219:LEU:HA	1.76	0.43
1:B:272:ARG:HD2	1:B:342:ASP:OD1	2.18	0.43
1:B:218:ASP:OD2	1:B:260:MET:CE	2.66	0.43
1:B:186:ARG:O	1:B:187:GLU:CD	2.56	0.43
1:A:87:ARG:CB	1:A:88:PRO:CD	2.74	0.43
1:A:126:LYS:HE3	1:A:126:LYS:HB2	1.76	0.43
1:B:303:ASP:CB	1:B:307:LYS:HD3	2.46	0.43
1:A:175:LEU:HB3	1:A:226:LEU:O	2.19	0.43
1:B:303:ASP:CG	1:B:307:LYS:HD2	2.39	0.43
1:B:99:LYS:N	1:B:386:PHE:CE2	2.83	0.43
1:A:283:MET:HG2	1:A:324:LEU:HD21	2.00	0.42
1:A:415:ALA:O	1:A:416:PRO:O	2.38	0.42
1:B:277:TRP:O	1:B:278:SER:C	2.56	0.42
1:A:156:VAL:O	1:A:156:VAL:HG13	2.19	0.42
1:A:417:SER:C	1:A:418:VAL:HG13	2.39	0.42
1:B:349:HIS:CG	1:B:350:PRO:CD	2.99	0.42
1:A:155:ILE:O	1:A:156:VAL:C	2.57	0.42
2:A:501:1SK:F31	2:A:501:1SK:CL28	2.63	0.42
1:A:204:MET:HE1	1:A:361:LEU:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:ARG:HD2	1:B:144:ARG:O	2.19	0.42
1:B:102:TYR:CE2	1:B:239:LEU:HD13	2.55	0.42
1:A:94:LEU:O	1:A:95:ARG:CG	2.65	0.42
1:A:135:LYS:O	1:A:139:HIS:HB2	2.20	0.42
1:B:125:LEU:HB2	1:B:168:LEU:HB2	2.01	0.42
1:B:416:PRO:O	1:B:417:SER:C	2.59	0.42
1:B:216:TYR:O	1:B:217:ARG:HB2	2.20	0.41
1:A:163:GLN:C	1:A:164:THR:HG23	2.39	0.41
1:A:277:TRP:CD1	1:A:277:TRP:C	2.93	0.41
1:B:102:TYR:CZ	1:B:239:LEU:HD13	2.56	0.41
1:B:319:GLU:N	1:B:319:GLU:CD	2.64	0.41
1:A:256:THR:OG1	1:A:258:GLU:OE1	2.39	0.41
1:A:119:ILE:CG1	1:A:376:GLU:HG2	2.43	0.41
1:A:159:ILE:HG22	1:A:160:TYR:CD1	2.56	0.41
1:A:271:ASN:OD1	1:A:271:ASN:C	2.59	0.41
1:A:349:HIS:CG	1:A:350:PRO:HD2	2.54	0.41
1:A:258:GLU:OE2	1:A:294:THR:CG2	2.68	0.41
1:A:330:ARG:NH1	1:A:330:ARG:CG	2.66	0.41
1:B:111:VAL:O	1:B:111:VAL:CG1	2.69	0.41
1:B:138:ALA:O	1:B:141:LYS:HB2	2.21	0.41
1:B:225:MET:CG	1:B:235:THR:HG21	2.47	0.41
1:B:271:ASN:C	1:B:273:ALA:H	2.24	0.41
1:B:409:LEU:CD1	1:B:410:GLY:N	2.66	0.41
1:A:214:ILE:HG22	1:A:215:ILE:N	2.36	0.40
1:A:323:LEU:HD12	1:A:323:LEU:HA	1.84	0.40
1:B:407:VAL:O	1:B:407:VAL:CG2	2.68	0.40
1:A:144:ARG:HD2	1:A:144:ARG:C	2.41	0.40
1:B:370:LYS:HA	1:B:371:PRO:HD2	1.91	0.40
1:A:127:LYS:CG	1:A:166:GLY:C	2.82	0.40
1:B:94:LEU:HB2	1:B:107:GLN:O	2.21	0.40
1:B:304:LYS:O	1:B:305:ILE:C	2.59	0.40
1:B:382:PHE:N	1:B:382:PHE:CD1	2.88	0.40
1:A:282:LEU:C	1:A:282:LEU:CD2	2.89	0.40
1:A:383:ASP:O	1:A:386:PHE:N	2.34	0.40
1:B:111:VAL:O	1:B:116:THR:HG21	2.21	0.40
1:B:131:VAL:CA	1:B:137:THR:CG2	2.82	0.40
1:A:282:LEU:HD22	1:A:282:LEU:C	2.42	0.40
1:B:136:ASP:O	1:B:137:THR:C	2.59	0.40
1:B:183:GLN:OE1	1:B:186:ARG:NH2	2.55	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	306/340 (90%)	276 (90%)	24 (8%)	6 (2%)	7	13
1	B	303/340 (89%)	269 (89%)	32 (11%)	2 (1%)	22	40
All	All	609/680 (90%)	545 (90%)	56 (9%)	8 (1%)	12	21

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	416	PRO
1	A	165	GLY
1	A	384	SER
1	B	417	SER
1	A	113	GLY
1	B	272	ARG
1	A	295	GLY
1	A	410	GLY

### 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	265/288 (92%)	240 (91%)	25 (9%)	8	16
1	B	262/288 (91%)	233 (89%)	29 (11%)	6	10
All	All	527/576 (92%)	473 (90%)	54 (10%)	8	13

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	87	ARG
1	A	108	VAL
1	A	126	LYS
1	A	132	ARG
1	A	232	VAL
1	A	242	GLU
1	A	257	ILE
1	A	258	GLU
1	A	282	LEU
1	A	294	THR
1	A	297	ASN
1	A	307	LYS
1	A	311	ASN
1	A	321	ARG
1	A	330	ARG
1	A	353	ARG
1	A	358[A]	GLU
1	A	358[B]	GLU
1	A	363	ARG
1	A	366	GLU
1	A	370	LYS
1	A	377	GLU
1	A	384	SER
1	A	392	VAL
1	A	406	GLN
1	B	107	GLN
1	B	112	THR
1	B	129	MET
1	B	130	ILE
1	B	131	VAL
1	B	133	ASN
1	B	140	THR
1	B	193[A]	ASP
1	B	193[B]	ASP
1	B	256	THR
1	B	258	GLU
1	B	264	ILE
1	B	267	ARG
1	B	278	SER
1	B	282	LEU
1	B	294	THR
1	B	297	ASN
1	B	299	LYS

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Mol	Chain	Res	Type
1	B	307	LYS
1	B	318	GLN
1	B	358	GLU
1	B	371	PRO
1	B	379	VAL
1	B	381	GLN
1	B	384	SER
1	B	392	VAL
1	B	393	ASP
1	B	405	ASN
1	B	418	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	ASN
1	A	297	ASN
1	A	311	ASN
1	A	406	GLN
1	B	208	HIS
1	B	228	HIS
1	B	297	ASN
1	B	381	GLN
1	B	405	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues 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
1	SEP	A	394	1	8,9,10	1.55	1 (12%)	8,12,14	2.10	1 (12%)
1	TPO	A	390	1	8,10,11	0.72	0	10,14,16	1.10	0
1	TPO	B	390	1	8,10,11	0.76	0	10,14,16	1.10	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
1	SEP	A	394	1	-	2/5/8/10	-
1	TPO	A	390	1	-	0/9/11/13	-
1	TPO	B	390	1	-	0/9/11/13	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	394	SEP	P-O1P	3.37	1.61	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	394	SEP	OG-CB-CA	5.28	113.28	108.14

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	394	SEP	N-CA-CB-OG
1	A	394	SEP	CA-CB-OG-P

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	394	SEP	3	0

## 5.5 Carbohydrates

There are no monosaccharides in this entry.

## 5.6 Ligand geometry

2 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	1SK	A	501	-	31,34,34	1.30	4 (12%)	42,50,50	2.61	14 (33%)
2	1SK	B	501	-	31,34,34	1.50	4 (12%)	42,50,50	2.85	20 (47%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	1SK	A	501	-	-	3/20/28/28	0/4/4/4
2	1SK	B	501	-	-	6/20/28/28	0/4/4/4

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	1SK	C4-C5	-5.39	1.38	1.44
2	A	501	1SK	C4-C5	-4.34	1.39	1.44
2	B	501	1SK	C7-C6	-2.91	1.39	1.43
2	B	501	1SK	C16-C15	-2.85	1.47	1.52
2	A	501	1SK	C19-CL28	2.55	1.79	1.73
2	B	501	1SK	C19-CL28	2.41	1.79	1.73
2	A	501	1SK	C7-C6	-2.25	1.40	1.43
2	A	501	1SK	C16-C15	-2.22	1.48	1.52

All (34) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	1SK	C2-N3-C6	7.98	121.02	114.81
2	A	501	1SK	C2-N3-C6	6.83	120.13	114.81
2	B	501	1SK	C4-C5-C6	6.57	120.02	115.87
2	A	501	1SK	C27-C18-C19	-6.56	117.60	121.88
2	B	501	1SK	C16-C15-N14	-5.80	104.72	113.44
2	A	501	1SK	C4-C5-C6	5.34	119.24	115.87
2	A	501	1SK	C18-C19-CL28	-4.95	115.98	121.68
2	B	501	1SK	N3-C2-N1	-4.56	121.55	128.68
2	B	501	1SK	C7-C11-N13	4.51	125.16	118.28
2	A	501	1SK	C16-C15-N14	-4.47	106.73	113.44
2	B	501	1SK	C10-C5-C4	-4.43	117.99	124.06
2	B	501	1SK	C7-C6-N3	4.08	121.84	118.75
2	B	501	1SK	O12-C11-C7	-4.02	115.48	120.22
2	A	501	1SK	N3-C2-N1	-4.02	122.40	128.68
2	A	501	1SK	C2-N1-C4	3.84	119.89	116.59
2	B	501	1SK	C27-C18-C19	-3.56	119.56	121.88
2	A	501	1SK	C15-C22-N23	-3.31	104.12	112.37
2	B	501	1SK	C2-N1-C4	3.30	119.42	116.59
2	A	501	1SK	F31-C27-C18	-3.30	106.96	112.70
2	A	501	1SK	C7-C6-N3	3.03	121.04	118.75
2	B	501	1SK	C18-C19-CL28	-3.00	118.23	121.68
2	A	501	1SK	C17-C18-C27	2.95	123.75	116.50
2	B	501	1SK	N14-C4-N1	2.76	122.25	118.06
2	A	501	1SK	C10-C5-C4	-2.70	120.36	124.06
2	B	501	1SK	F30-C27-C18	-2.65	108.09	112.70
2	B	501	1SK	C5-C4-N14	-2.59	118.19	120.63
2	B	501	1SK	C15-C22-N23	-2.35	106.52	112.37
2	B	501	1SK	O12-C11-N13	-2.30	119.31	122.58
2	B	501	1SK	C5-C6-N3	-2.18	118.42	122.54
2	B	501	1SK	C5-C4-N1	-2.17	119.56	121.35
2	B	501	1SK	C20-C21-C16	-2.12	119.06	121.20
2	B	501	1SK	C17-C18-C27	2.10	121.66	116.50
2	A	501	1SK	C5-C6-N3	-2.05	118.67	122.54
2	A	501	1SK	C5-C4-N1	-2.04	119.67	121.35

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	1SK	C16-C15-N14-C4
2	B	501	1SK	C16-C15-C22-N23
2	B	501	1SK	C22-C15-N14-C4
2	B	501	1SK	C16-C15-N14-C4

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Mol	Chain	Res	Type	Atoms
2	B	501	1SK	C5-C4-N14-C15
2	A	501	1SK	N14-C15-C22-N23
2	B	501	1SK	N14-C15-C22-N23
2	A	501	1SK	C22-C15-N14-C4
2	B	501	1SK	N1-C4-N14-C15

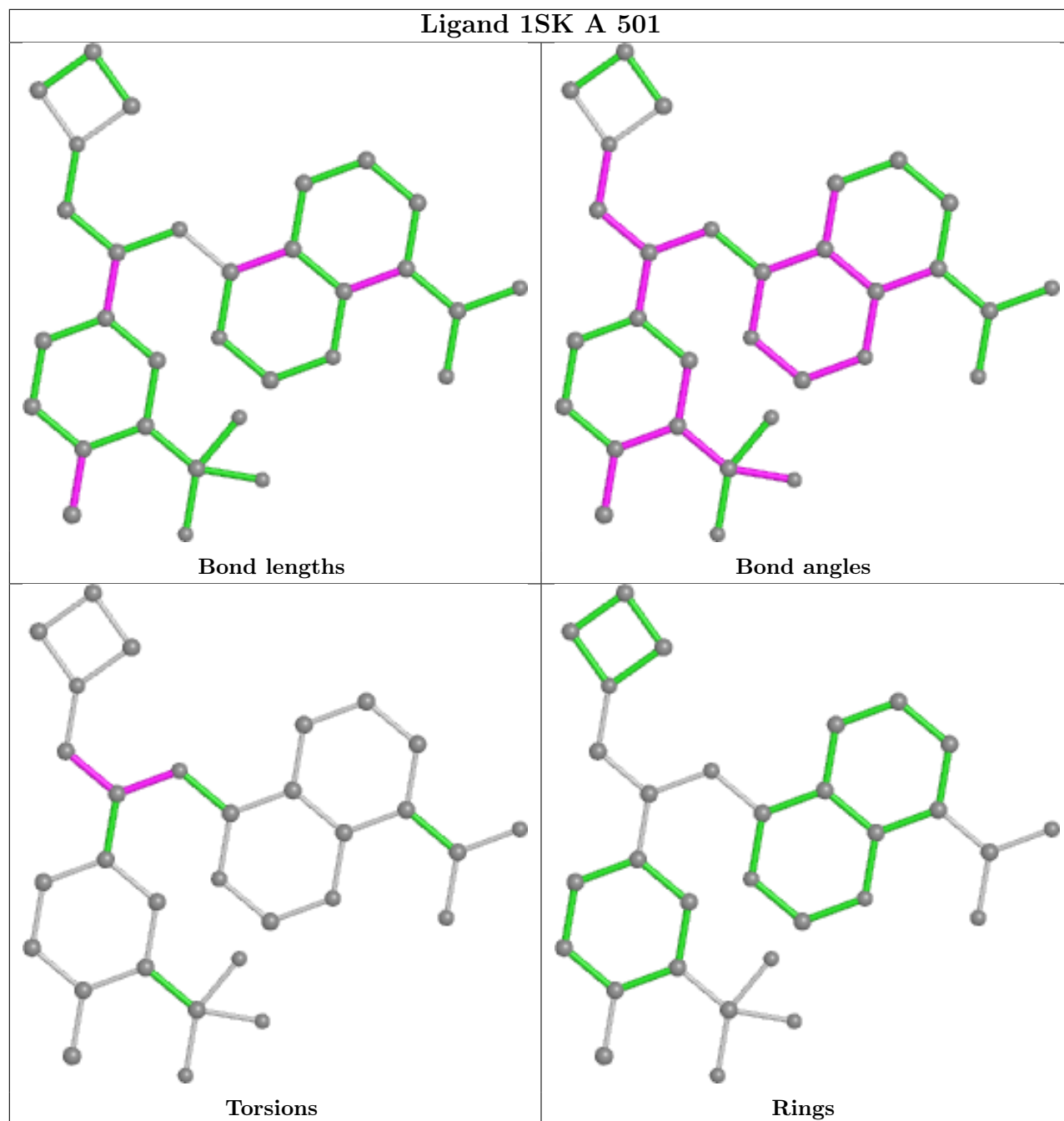
There are no ring outliers.

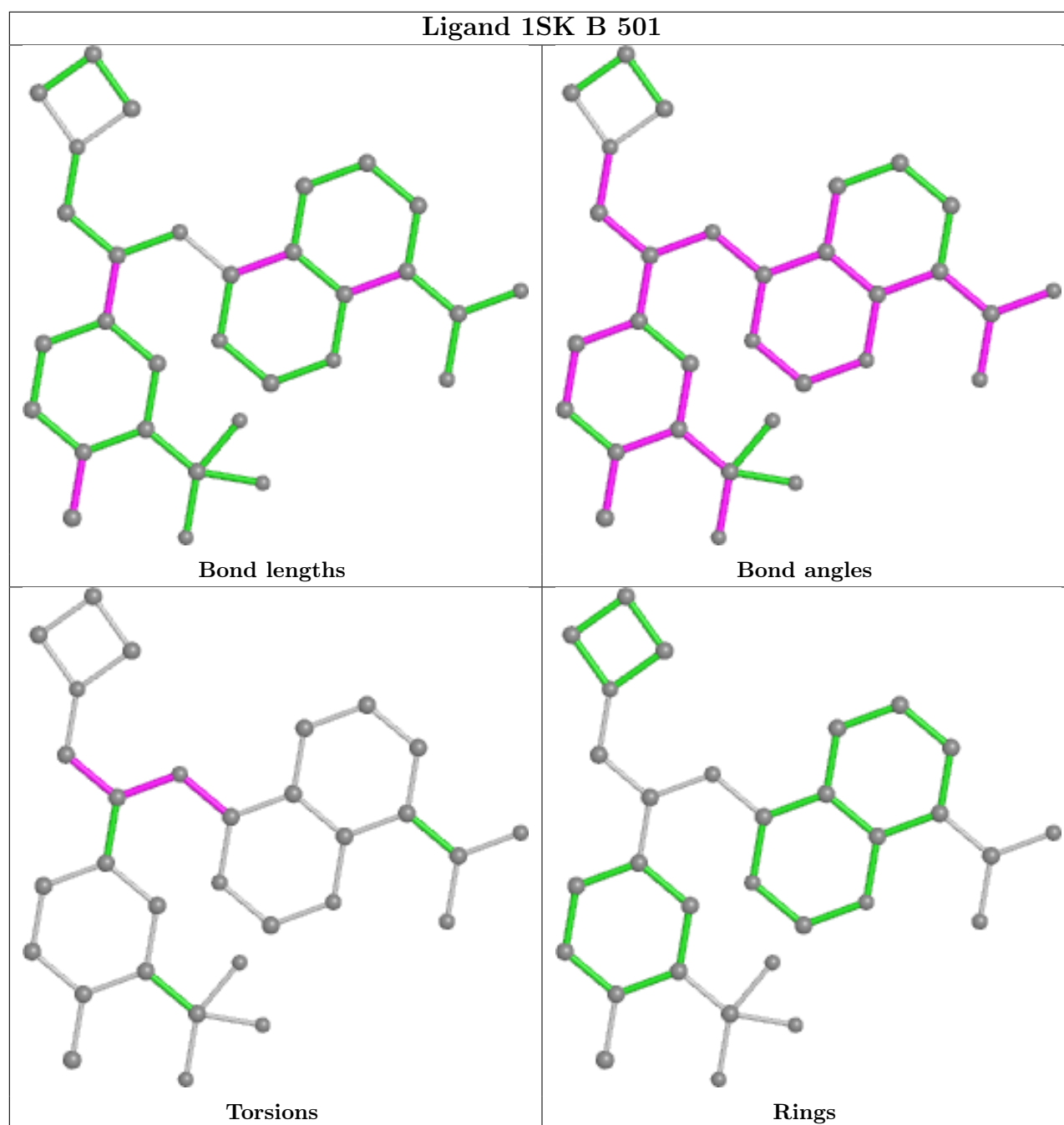
2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	1SK	3	0
2	B	501	1SK	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

## Ligand 1SK A 501





## 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, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	310/340 (91%)	0.08	4 (1%) 77 82	35, 49, 64, 76	0
1	B	308/340 (90%)	0.31	18 (5%) 23 26	36, 50, 65, 71	0
All	All	618/680 (90%)	0.19	22 (3%) 42 47	35, 49, 65, 76	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	85	LYS	6.7
1	B	86	ILE	6.3
1	B	102	TYR	3.9
1	B	168	LEU	3.8
1	B	132	ARG	3.4
1	B	308	CYS	3.3
1	B	128	ALA	3.2
1	A	417	SER	3.1
1	B	113	GLY	3.0
1	B	328	LEU	2.7
1	A	388	ARG	2.5
1	B	87	ARG	2.4
1	B	166	GLY	2.4
1	B	127	LYS	2.4
1	B	264	ILE	2.4
1	A	405	ASN	2.4
1	B	129	MET	2.3
1	B	126	LYS	2.3
1	A	369	PHE	2.3
1	B	267	ARG	2.1
1	B	130	ILE	2.0
1	B	310	LEU	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
1	SEP	A	394	10/11	0.66	0.21	59,60,61,61	4
1	TPO	A	390	11/12	0.80	0.21	61,61,61,62	4
1	TPO	B	390	11/12	0.86	0.16	62,63,64,64	4

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

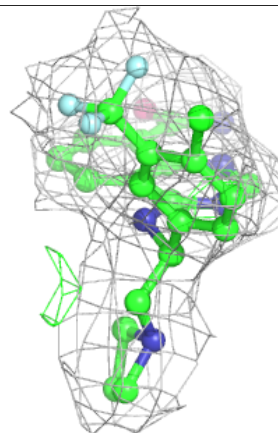
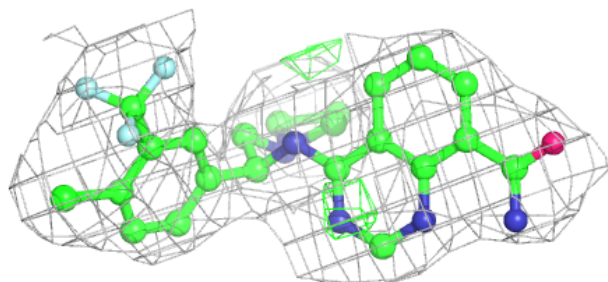
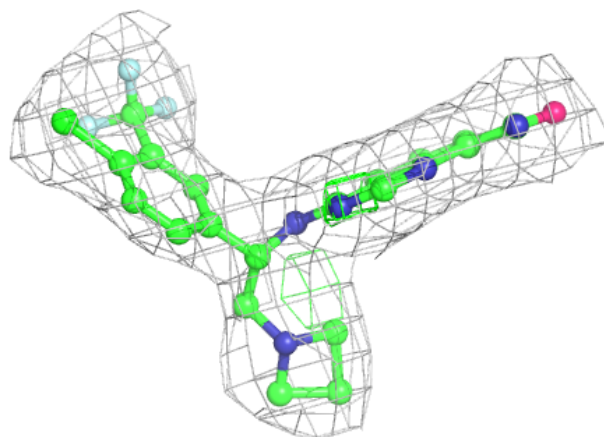
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	1SK	A	501	31/31	0.90	0.18	48,59,68,71	0
2	1SK	B	501	31/31	0.92	0.23	45,67,77,78	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.

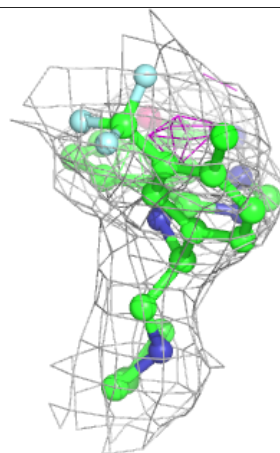
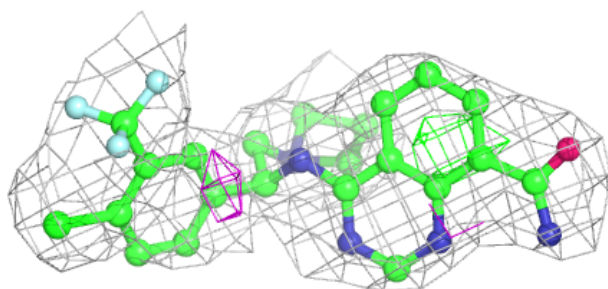
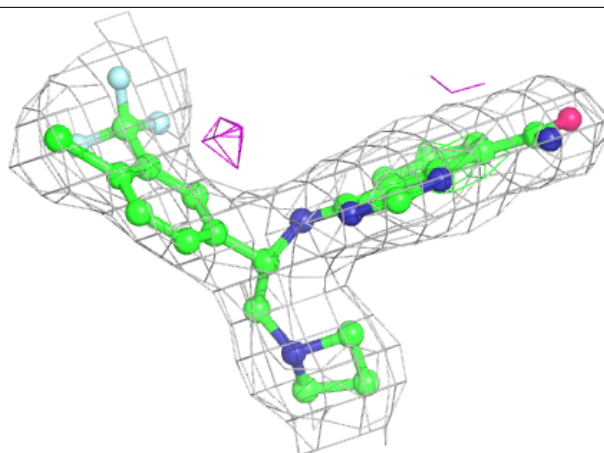
**Electron density around 1SK A 501:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around 1SK B 501:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.