



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:03 PM GMT

PDB ID : 3SRK  
Title : A new class of suicide inhibitor blocks nucleotide binding to pyruvate kinase  
Authors : Morgan, H.P.; Walsh, M.; Blackburn, E.A.; Wear, M.A.; Boxer, M.; Shen, M.; McNae, I.W.; Michels, P.A.M.; Auld, D.S.; Fothergill-Gilmore, L.A.; Walkinshaw, M.D.  
Deposited on : 2011-07-07  
Resolution : 2.65 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

---

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

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

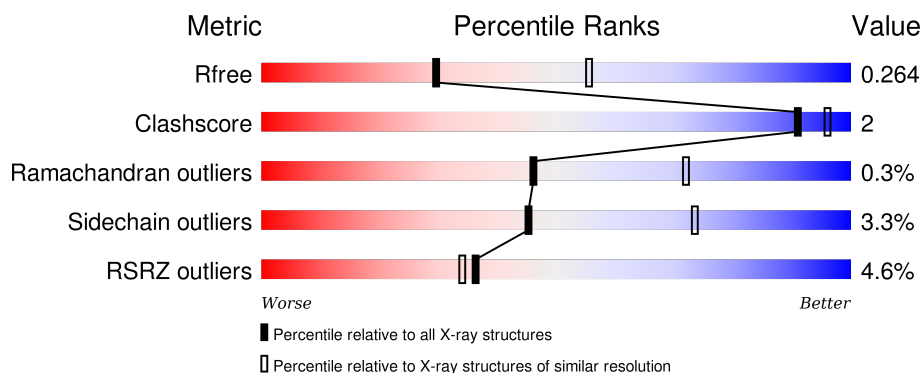
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

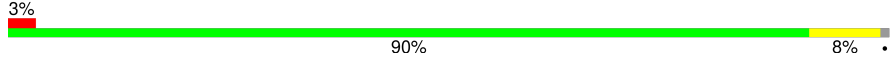
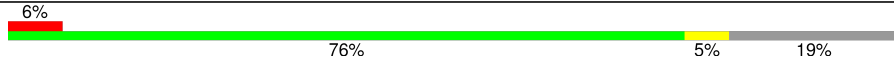
The reported resolution of this entry is 2.65 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	3152 (2.70-2.62)
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)
RSRZ outliers	91569	3161 (2.70-2.62)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	499	 3% 90% 8%
1	B	499	 6% 76% 5% 19%

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	LSA	A	500	-	-	-	X
2	LSA	B	500	-	-	-	X
4	PTK	A	502	-	-	-	X

## 2 Entry composition [i](#)

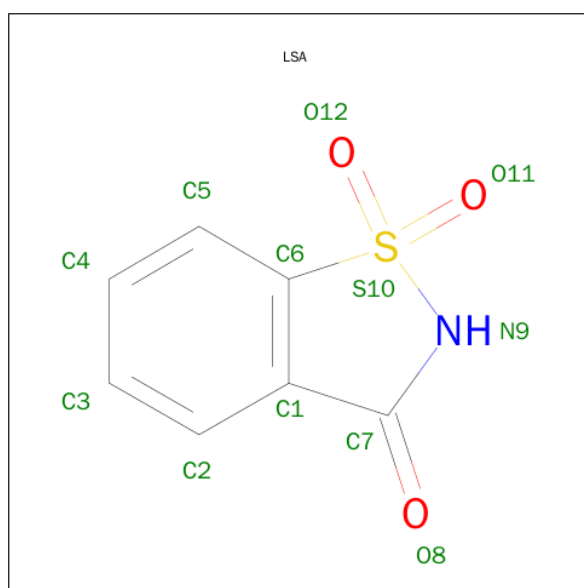
There are 5 unique types of molecules in this entry. The entry contains 7061 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Pyruvate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	493	Total	C	N	O	S	0	1	0
			3768	2349	663	729	27			
1	B	405	Total	C	N	O	S	0	3	0
			3098	1934	546	594	24			

- Molecule 2 is 1,2-BENZISOTHIAZOL-3(2H)-ONE 1,1-DIOXIDE (three-letter code: LSA) (formula: C<sub>7</sub>H<sub>5</sub>NO<sub>3</sub>S).

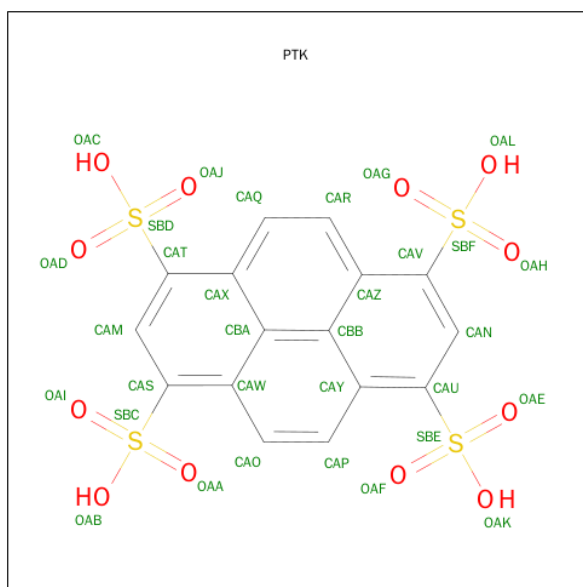


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			11	7	1	2	1		
2	B	1	Total	C	N	O	S	0	0
			11	7	1	2	1		

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	2	Total K 2 2	0	0
3	A	2	Total K 2 2	0	0

- Molecule 4 is PYRENE-1,3,6,8-TETRASULFONIC ACID (three-letter code: PTK) (formula: C<sub>16</sub>H<sub>10</sub>O<sub>12</sub>S<sub>4</sub>).



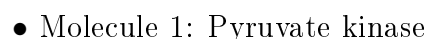
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O S 32 16 12 4	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	69	Total O 69 69	0	0
5	B	68	Total O 68 68	0	0



- Molecule 1: Pyruvate kinase



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	122.42Å 130.16Å 166.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.31 – 2.65 39.30 – 2.65	Depositor EDS
% Data completeness (in resolution range)	97.2 (39.31-2.65) 97.2 (39.30-2.65)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.84 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, $R_{free}$	0.223 , 0.266 0.225 , 0.264	Depositor DCC
$R_{free}$ test set	1901 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	55.5	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 48.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 37835 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	7061	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.32	0/3827	0.46	0/5179
1	B	0.31	0/3145	0.50	3/4253 (0.1%)
All	All	0.32	0/6972	0.48	3/9432 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	180	PRO	CB-CA-C	-9.96	87.10	112.00
1	B	180	PRO	N-CA-C	5.84	127.29	112.10
1	B	1	SER	CB-CA-C	5.81	121.13	110.10

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	3768	0	3766	15	0
1	B	3098	0	3113	8	0
2	A	11	0	5	0	0
2	B	11	0	5	1	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	32	0	8	8	0
5	A	69	0	0	0	0
5	B	68	0	0	0	0
All	All	7061	0	6897	31	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 2.

All (31) 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)
4:A:502:PTK:OAF	4:A:502:PTK:HAP	1.47	1.07
4:A:502:PTK:OAH	4:A:502:PTK:HAR	1.47	1.06
4:A:502:PTK:OAI	4:A:502:PTK:HAO	1.62	0.95
4:A:502:PTK:OAJ	4:A:502:PTK:HAQ	1.61	0.95
4:A:502:PTK:OAH	4:A:502:PTK:CAR	2.30	0.76
4:A:502:PTK:OAF	4:A:502:PTK:CAP	2.30	0.75
4:A:502:PTK:OAJ	4:A:502:PTK:CAQ	2.41	0.67
4:A:502:PTK:OAI	4:A:502:PTK:CAO	2.41	0.67
1:A:89:ILE:HG23	1:A:128:TYR:HB2	1.79	0.65
1:B:298:MET:HE3	1:B:316:VAL:HG22	1.86	0.57
1:A:338:TYR:HB3	1:A:341:GLU:HB2	1.86	0.57
1:B:0:MET:HB3	1:B:360:TYR:HD2	1.71	0.56
1:B:394:LYS:HB2	1:B:470:VAL:HG12	1.88	0.55
1:A:24:ILE:HB	1:A:328:MET:HG3	1.90	0.54
1:A:142:TYR:HB3	1:A:146:GLY:HA2	1.92	0.51
1:A:481:ALA:HA	1:A:489:ALA:HA	1.91	0.51
1:B:24:ILE:HG12	1:B:47:VAL:HB	1.97	0.47
1:B:304[B]:TYR:N	1:B:304[B]:TYR:CD2	2.83	0.46
1:A:99:ALA:HB3	1:A:171:ILE:HG13	1.97	0.46
1:A:148:LEU:HB2	1:A:169:HIS:HD2	1.81	0.46
1:A:17:ASN:H	1:A:17:ASN:ND2	2.16	0.44
1:B:51:ASN:HB3	2:B:500:LSA:O12	2.18	0.44
1:B:301:SER:HB3	1:B:312:GLU:OE1	2.18	0.43
1:B:47:VAL:HG22	1:B:79:ALA:HB3	2.00	0.43
1:A:179:LEU:HB3	1:A:182:CYS:HB2	2.01	0.43
1:A:17:ASN:H	1:A:17:ASN:HD22	1.67	0.42
1:A:112:ASP:HA	1:A:113:PRO:HD3	1.87	0.42
1:A:47:VAL:HG22	1:A:79:ALA:HB3	2.02	0.42
1:A:79:ALA:HB2	1:A:429:ARG:O	2.19	0.42
1:A:398:VAL:HG22	1:A:479:ILE:HG23	2.01	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:ARG:H	1:A:139:ASN:ND2	2.19	0.41

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	490/499 (98%)	475 (97%)	13 (3%)	2 (0%)	39	65
1	B	402/499 (81%)	392 (98%)	9 (2%)	1 (0%)	52	77
All	All	892/998 (89%)	867 (97%)	22 (2%)	3 (0%)	46	72

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	296	THR
1	B	296	THR
1	A	44	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	413/417 (99%)	397 (96%)	16 (4%)	39	67
1	B	339/417 (81%)	330 (97%)	9 (3%)	52	80

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	752/834 (90%)	727 (97%)	25 (3%)	45 73

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	129	GLN
1	A	136	ARG
1	A	173	ASP
1	A	175	ARG
1	A	177	VAL
1	A	194	ARG
1	A	262	ARG
1	A	390	GLU
1	A	426	GLN
1	A	435	GLN
1	A	453	LYS
1	A	471	GLN
1	A	479	ILE
1	A	488	TYR
1	A	490	ASN
1	A	492	THR
1	B	0	MET
1	B	17	ASN
1	B	34	VAL
1	B	40	LEU
1	B	224	LYS
1	B	262	ARG
1	B	446	LYS
1	B	454	GLU
1	B	479	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	139	ASN
1	A	169	HIS
1	A	198	GLN
1	A	203	GLN
1	A	242	HIS
1	A	305	ASN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	386	ASN
1	A	435	GLN
1	A	471	GLN
1	A	490	ASN
1	B	17	ASN
1	B	370	GLN
1	B	386	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	$\# Z  > 2$
2	LSA	A	500	1	11,12,13	4.00	5 (45%)	16,18,20	2.97	5 (31%)
4	PTK	A	502	-	35,35,35	2.54	17 (48%)	50,60,60	1.89	12 (24%)
2	LSA	B	500	1	11,12,13	3.84	5 (45%)	16,18,20	3.27	5 (31%)

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	LSA	A	500	1	-	0/0/12/15	0/2/2/2
4	PTK	A	502	-	-	0/24/24/24	0/4/4/4
2	LSA	B	500	1	-	0/0/12/15	0/2/2/2

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	LSA	C7-N9	-7.73	1.34	1.45
2	B	500	LSA	C7-N9	-7.52	1.35	1.45
4	A	502	PTK	CAU-SBE	-5.21	1.70	1.78
4	A	502	PTK	CAV-SBF	-5.21	1.70	1.78
4	A	502	PTK	CAS-SBC	-4.54	1.71	1.78
4	A	502	PTK	CAU-CAY	-4.52	1.36	1.43
4	A	502	PTK	CAV-CAZ	-4.45	1.36	1.43
4	A	502	PTK	CAT-SBD	-4.15	1.72	1.78
4	A	502	PTK	CAT-CAX	-4.12	1.37	1.43
4	A	502	PTK	CAS-CAW	-4.06	1.37	1.43
2	B	500	LSA	C6-S10	-2.98	1.72	1.75
2	B	500	LSA	S10-N9	-2.87	1.61	1.63
4	A	502	PTK	CAW-CBA	-2.44	1.36	1.42
4	A	502	PTK	CAY-CBB	-2.42	1.37	1.42
4	A	502	PTK	CAX-CBA	-2.40	1.37	1.42
2	A	500	LSA	S10-N9	-2.40	1.61	1.63
4	A	502	PTK	CAZ-CBB	-2.36	1.37	1.42
2	A	500	LSA	C6-S10	-2.27	1.73	1.75
4	A	502	PTK	CBA-CBB	-2.26	1.36	1.43
4	A	502	PTK	CAR-CAZ	-2.23	1.37	1.42
4	A	502	PTK	CAP-CAY	-2.18	1.37	1.42
4	A	502	PTK	CAQ-CAX	-2.06	1.38	1.42
4	A	502	PTK	CAO-CAW	-2.05	1.38	1.42
2	B	500	LSA	O12-S10	6.35	1.50	1.43
2	B	500	LSA	O11-S10	6.56	1.50	1.43
2	A	500	LSA	O12-S10	7.00	1.51	1.43
2	A	500	LSA	O11-S10	7.15	1.51	1.43

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	LSA	C1-C6-S10	-9.80	103.50	110.32
2	A	500	LSA	C1-C6-S10	-9.26	103.87	110.32

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	LSA	O11-S10-O12	-4.89	109.40	116.22
4	A	502	PTK	CAP-CAY-CAU	-3.61	119.04	123.83
2	A	500	LSA	O11-S10-O12	-3.58	111.23	116.22
4	A	502	PTK	CAR-CAZ-CAV	-3.57	119.08	123.83
4	A	502	PTK	CAO-CAW-CAS	-3.46	119.24	123.83
4	A	502	PTK	CAQ-CAX-CAT	-3.42	119.29	123.83
2	A	500	LSA	C7-C1-C2	2.07	132.79	128.84
4	A	502	PTK	OAJ-SBD-CAT	2.19	108.68	106.20
4	A	502	PTK	CAN-CAU-SBE	2.19	121.62	117.82
4	A	502	PTK	CAN-CAV-SBF	2.19	121.62	117.82
2	A	500	LSA	O12-S10-N9	2.44	113.50	110.98
4	A	502	PTK	OAI-SBC-CAS	2.71	109.27	106.20
2	B	500	LSA	O12-S10-N9	2.85	113.92	110.98
4	A	502	PTK	OAA-SBC-CAS	2.92	109.51	106.20
2	B	500	LSA	C6-S10-N9	2.96	97.39	94.38
4	A	502	PTK	OAD-SBD-CAT	3.37	110.02	106.20
4	A	502	PTK	OAE-SBE-CAU	4.63	111.45	106.20
4	A	502	PTK	OAG-SBF-CAV	4.63	111.45	106.20
2	A	500	LSA	C5-C6-S10	4.72	134.49	127.12
2	B	500	LSA	C5-C6-S10	5.00	134.93	127.12

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	502	PTK	8	0
2	B	500	LSA	1	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, 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	493/499 (98%)	0.07	13 (2%) 59 58	37, 60, 92, 105	0
1	B	405/499 (81%)	0.36	28 (6%) 20 17	38, 60, 94, 156	0
All	All	898/998 (89%)	0.20	41 (4%) 36 34	37, 60, 94, 156	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	179	LEU	6.8
1	B	182	CYS	6.4
1	B	180	PRO	6.0
1	B	449	HIS	5.6
1	B	181	GLY	5.5
1	B	447	LEU	5.4
1	B	177	VAL	5.0
1	B	178	ASN	4.6
1	B	452	GLY	4.2
1	B	488	TYR	4.0
1	B	445	ASP	3.8
1	B	184	VAL	3.7
1	A	168	SER	3.6
1	B	444	ALA	3.5
1	A	488	TYR	3.2
1	B	371[A]	HIS	3.2
1	B	183	ASP	3.1
1	B	442	PHE	3.1
1	B	88	GLU	3.0
1	B	451	GLU	3.0
1	A	447	LEU	2.9
1	A	480	HIS	2.9
1	A	372	ILE	2.8
1	B	443	ASP	2.7

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	155	HIS	2.7
1	B	372	ILE	2.7
1	B	446	LYS	2.7
1	B	187	PRO	2.6
1	A	158	GLU	2.6
1	B	320	VAL	2.5
1	B	448	GLY	2.5
1	B	185	ASP	2.4
1	A	153	GLN	2.4
1	B	186	LEU	2.4
1	B	317	ALA	2.3
1	B	319	ALA	2.3
1	A	100	VAL	2.2
1	A	103	ARG	2.1
1	A	449[A]	HIS	2.1
1	A	307	ARG	2.0
1	A	481	ALA	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	LSA	B	500	11/12	0.91	0.41	9.39	41,42,43,43	0
2	LSA	A	500	11/12	0.90	0.51	9.14	40,41,42,42	0
4	PTK	A	502	32/32	0.79	0.30	3.76	63,64,66,66	32
3	K	B	499	1/1	0.98	0.09	-3.88	48,48,48,48	0

*Continued on next page...*



*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	K	A	501	1/1	0.97	0.04	-6.17	51,51,51,51	0
3	K	A	499	1/1	0.94	0.14	-	72,72,72,72	0
3	K	B	501	1/1	0.93	0.07	-	76,76,76,76	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.