



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:57 PM GMT

PDB ID : 4B2Z
Title : Structure of Osh6 in complex with phosphatidylserine
Authors : Maeda, K.; Anand, K.; Chiapparino, A.; Kumar, A.; Poletto, M.; Kaksonen, M.; Gavin, A.C.
Deposited on : 2012-07-19
Resolution : 1.95 Å(reported)

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

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

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

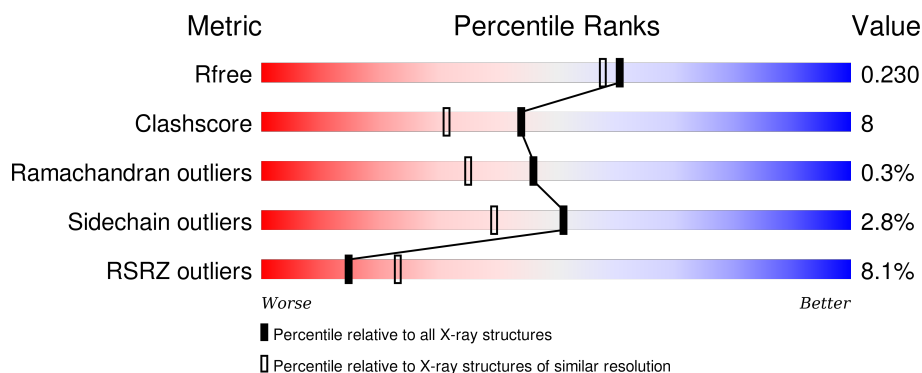
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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

Mol	Chain	Length	Quality of chain
1	A	448	<div> <div>9%</div> <div>77%</div> <div>10%</div> <div>11%</div> </div>
1	B	448	<div> <div>6%</div> <div>78%</div> <div>10%</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
2	SO4	A	1435	-	-	-	X
4	DTU	A	1437	-	-	-	X
5	DTT	B	1437	-	-	X	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7029 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 OXYSTEROL-BINDING PROTEIN HOMOLOG 6.

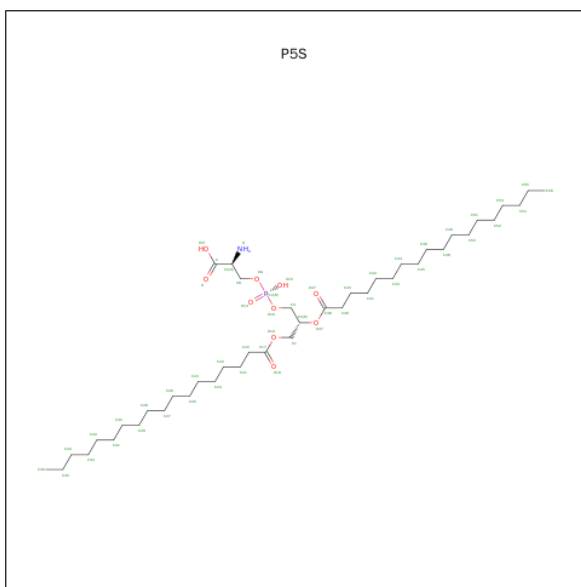
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	399	Total	C	N	O	S	15	0	0
			3264	2102	548	603	11			
1	B	400	Total	C	N	O	S	7	0	0
			3269	2105	549	604	11			

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



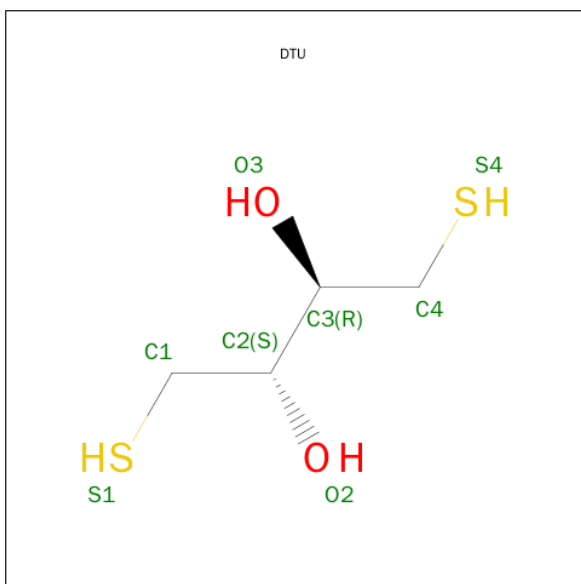
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is O-[(R)-{[(2R)-2,3-BIS(OCTADECANOYLOXY)PROPYL]OXY}(HYDROXY)PHOSPHORYL]-L-SERINE (three-letter code: P5S) (formula: C₄₂H₈₂NO₁₀P).



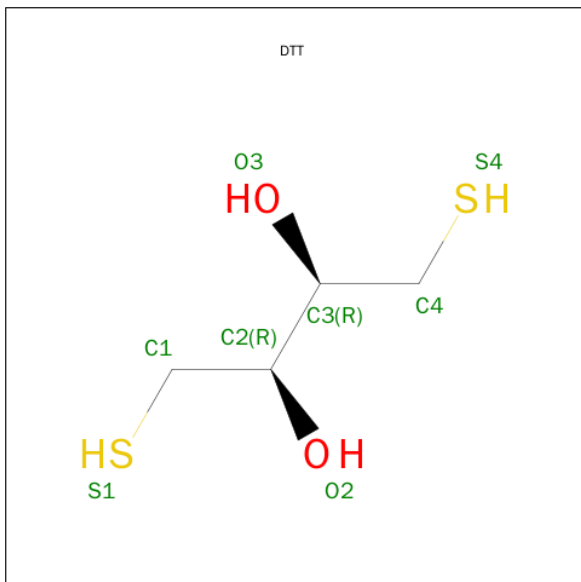
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			53	41	1	10	1		
3	B	1	Total	C	N	O	P	0	0
			54	42	1	10	1		

- Molecule 4 is (2R,3S)-1,4-DIMERCAPTOBUTANE-2,3-DIOL (three-letter code: DTU) (formula: $C_4H_{10}O_2S_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	S	0	0
			8	4	2	2		

- Molecule 5 is 2,3-DIHYDROXY-1,4-DITHIOBUTANE (three-letter code: DTT) (formula: $C_4H_{10}O_2S_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	O	S	0	0
			8	4	2	2		

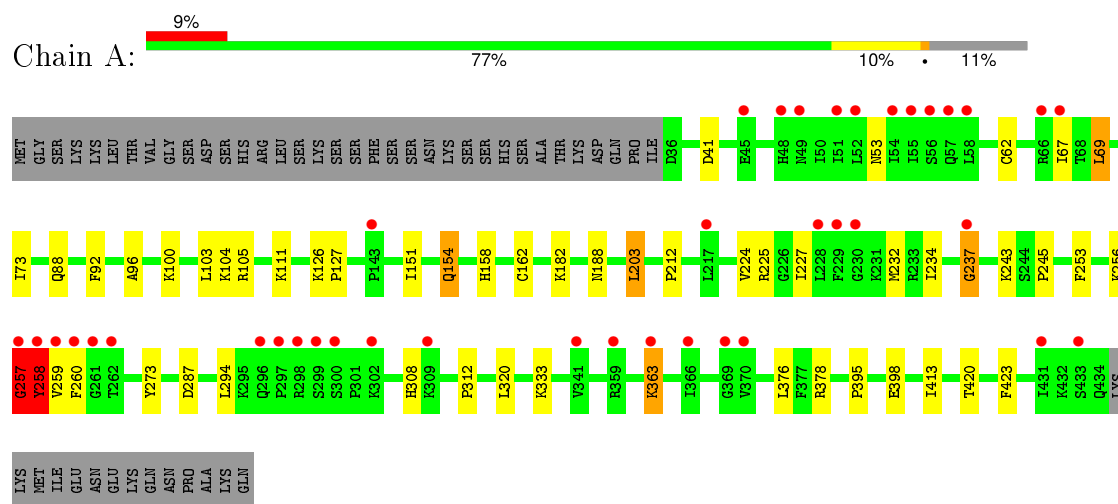
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	167	Total	O	0	0
			167	167		
6	B	196	Total	O	0	0
			196	196		

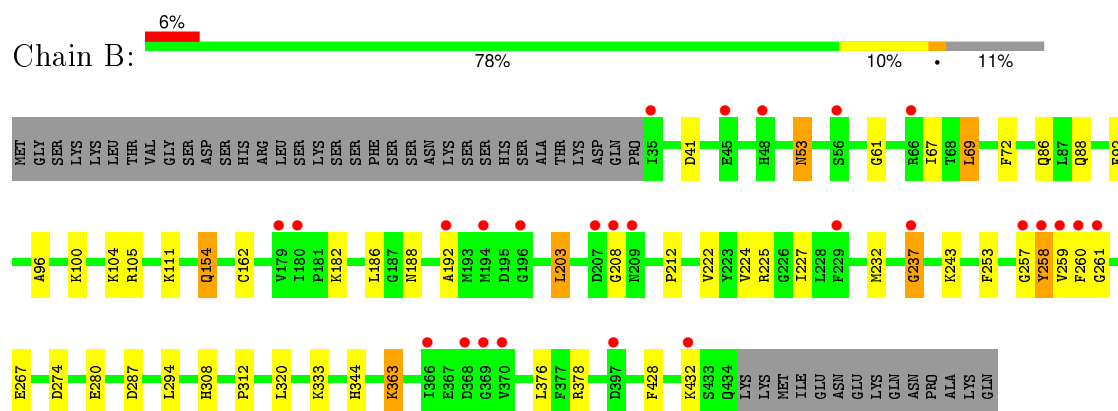
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: OXYSTEROL-BINDING PROTEIN HOMOLOG 6



• Molecule 1: OXYSTEROL-BINDING PROTEIN HOMOLOG 6



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.00Å 72.60Å 122.90Å 90.00° 111.40° 90.00°	Depositor
Resolution (Å)	46.57 – 1.95 46.57 – 1.95	Depositor EDS
% Data completeness (in resolution range)	91.8 (46.57-1.95) 91.8 (46.57-1.95)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 1.95Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.198 , 0.236 0.194 , 0.230	Depositor DCC
R_{free} test set	1318 reflections (2.04%)	DCC
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.462	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 65896 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7029	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: DTT, SO4, DTU, P5S

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	0.39	0/3350	0.56	3/4531 (0.1%)
1	B	0.40	0/3355	0.56	1/4538 (0.0%)
All	All	0.40	0/6705	0.56	4/9069 (0.0%)

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	A	0	1
1	B	0	2
All	All	0	3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	259	VAL	CB-CA-C	-6.63	98.80	111.40
1	A	258	TYR	N-CA-C	6.07	127.38	111.00
1	A	237	GLY	N-CA-C	5.18	126.05	113.10
1	B	237	GLY	N-CA-C	5.03	125.68	113.10

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	257	GLY	Peptide
1	B	257	GLY	Peptide

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Mol	Chain	Res	Type	Group
1	B	260	PHE	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	3264	0	3247	51	0
1	B	3269	0	3249	50	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	53	0	75	8	0
3	B	54	0	80	12	0
4	A	8	0	10	3	0
5	B	8	0	10	5	0
6	A	167	0	0	4	0
6	B	196	0	0	7	1
All	All	7029	0	6671	106	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 8.

All (106) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:GLY:HA2	5:B:1437:DTT:S1	2.14	0.88
1:A:227:ILE:HG22	1:A:232:MET:HE2	1.59	0.85
1:A:245:PRO:HA	5:B:1437:DTT:H12	1.61	0.81
1:A:227:ILE:HG22	1:A:232:MET:CE	2.11	0.80
1:B:227:ILE:HG22	1:B:232:MET:CE	2.11	0.80
1:A:182:LYS:HE2	6:B:2182:HOH:O	1.81	0.80
6:A:2153:HOH:O	1:B:182:LYS:HE2	1.82	0.80
1:B:53:ASN:HB3	6:B:2011:HOH:O	1.82	0.79
1:B:227:ILE:HG22	1:B:232:MET:HE2	1.68	0.74
1:B:267:GLU:HG3	6:B:2124:HOH:O	1.90	0.71
1:B:237:GLY:O	6:B:2113:HOH:O	2.08	0.70
1:A:260:PHE:O	1:A:260:PHE:CD1	2.44	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:VAL:HG21	1:A:232:MET:HE1	1.73	0.69
1:A:363:LYS:HE3	1:A:363:LYS:HA	1.75	0.69
1:B:222:VAL:HG12	3:B:1436:P5S:H44	1.77	0.66
1:B:363:LYS:HE3	1:B:363:LYS:HA	1.78	0.65
1:B:92:PHE:CE1	1:B:111:LYS:HE2	2.34	0.63
3:B:1436:P5S:H41	3:B:1436:P5S:H45	1.82	0.62
1:A:127:PRO:HD2	3:A:1436:P5S:O18	1.99	0.61
1:B:237:GLY:O	1:B:253:PHE:O	2.19	0.61
1:A:258:TYR:H	1:A:258:TYR:HD1	1.45	0.61
1:A:188:ASN:O	1:A:227:ILE:HG12	2.01	0.61
1:A:420:THR:HG23	4:A:1437:DTU:H3	1.83	0.60
1:A:423:PHE:HB2	4:A:1437:DTU:O2	2.01	0.60
3:B:1436:P5S:H45	3:B:1436:P5S:C41	2.32	0.59
1:A:154:GLN:NE2	1:A:162:CYS:HB3	2.18	0.59
1:A:62:CYS:HB3	1:B:208:GLY:HA3	1.85	0.58
1:A:224:VAL:HG21	1:A:232:MET:CE	2.32	0.58
1:B:224:VAL:HG21	1:B:232:MET:CE	2.33	0.58
1:A:92:PHE:CE1	1:A:111:LYS:HE2	2.40	0.57
1:A:237:GLY:O	1:A:253:PHE:O	2.22	0.56
1:B:227:ILE:HG22	1:B:232:MET:HE3	1.85	0.55
1:B:224:VAL:HG21	1:B:232:MET:HE1	1.87	0.55
1:A:243:LYS:HD3	5:B:1437:DTT:S4	2.48	0.54
1:A:363:LYS:CA	1:A:363:LYS:HE3	2.37	0.53
1:B:188:ASN:O	1:B:227:ILE:HG12	2.08	0.53
1:B:222:VAL:O	3:B:1436:P5S:H43A	2.08	0.53
1:A:225:ARG:NH1	1:B:274:ASP:HA	2.24	0.52
1:B:428:PHE:O	1:B:432:LYS:HG3	2.09	0.52
1:A:234:ILE:HD11	3:A:1436:P5S:H45	1.93	0.51
1:B:154:GLN:NE2	1:B:162:CYS:HB3	2.25	0.51
1:A:258:TYR:N	1:A:258:TYR:CD1	2.78	0.51
1:A:287:ASP:OD1	1:A:308:HIS:CE1	2.64	0.51
1:A:420:THR:OG1	4:A:1437:DTU:H2	2.10	0.51
1:A:73:ILE:HB	3:A:1436:P5S:H49A	1.93	0.51
3:B:1436:P5S:O18	3:B:1436:P5S:H22A	2.10	0.50
1:A:227:ILE:HG22	1:A:232:MET:HE3	1.93	0.49
1:B:363:LYS:HE3	1:B:363:LYS:CA	2.42	0.49
1:B:258:TYR:N	1:B:258:TYR:CD1	2.75	0.49
1:A:287:ASP:OD1	1:A:308:HIS:HE1	1.95	0.49
1:A:88:GLN:HE21	1:A:312:PRO:HB2	1.78	0.48
1:B:237:GLY:HA2	6:B:2106:HOH:O	2.13	0.48
1:B:192:ALA:HB2	3:B:1436:P5S:H20	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:320:LEU:HD11	1:A:333:LYS:HE3	1.95	0.48
1:A:73:ILE:HB	3:A:1436:P5S:C49	2.44	0.47
1:A:126:LYS:HB3	3:A:1436:P5S:H48	1.95	0.47
1:A:260:PHE:O	1:A:260:PHE:HD1	1.95	0.47
1:B:320:LEU:HD11	1:B:333:LYS:HE3	1.96	0.47
1:B:287:ASP:OD1	1:B:308:HIS:CE1	2.67	0.47
1:A:256:LYS:C	1:A:257:GLY:O	2.53	0.47
1:A:273:TYR:OH	5:B:1437:DTT:H41	2.15	0.47
6:A:2092:HOH:O	1:B:243:LYS:NZ	2.48	0.47
1:A:256:LYS:HG2	1:A:257:GLY:O	2.15	0.47
1:B:192:ALA:CB	3:B:1436:P5S:H20	2.45	0.46
1:B:100:LYS:O	1:B:105:ARG:NH2	2.47	0.46
1:B:203:LEU:HA	1:B:212:PRO:HB3	1.97	0.46
1:B:287:ASP:OD1	1:B:308:HIS:HE1	1.98	0.46
3:B:1436:P5S:H24	3:B:1436:P5S:H21	1.62	0.46
1:A:203:LEU:HA	1:A:212:PRO:HB3	1.98	0.46
1:B:222:VAL:HG11	3:B:1436:P5S:H21A	1.97	0.45
1:B:86:GLN:NE2	6:B:2028:HOH:O	2.48	0.45
6:A:2101:HOH:O	1:B:243:LYS:HE2	2.15	0.45
1:B:88:GLN:HE21	1:B:312:PRO:HB2	1.80	0.45
1:B:104:LYS:HE2	1:B:104:LYS:HB3	1.72	0.45
1:B:222:VAL:CG1	3:B:1436:P5S:H44	2.45	0.45
1:B:378:ARG:HG3	1:B:378:ARG:HH11	1.82	0.45
1:A:395:PRO:HB2	1:A:398:GLU:HB2	1.98	0.45
1:B:280:GLU:HG3	6:B:2121:HOH:O	2.16	0.44
1:B:67:ILE:CD1	1:B:69:LEU:HD13	2.47	0.44
1:A:126:LYS:HG2	3:A:1436:P5S:H49	1.98	0.44
1:A:225:ARG:O	1:A:232:MET:HA	2.16	0.44
1:A:158:HIS:HD2	6:A:2071:HOH:O	2.00	0.44
1:A:100:LYS:O	1:A:105:ARG:NH2	2.50	0.44
1:A:151:ILE:HD11	1:A:413:ILE:HD11	2.00	0.44
1:A:67:ILE:O	1:A:67:ILE:HD12	2.18	0.43
1:A:67:ILE:CD1	1:A:69:LEU:HD13	2.48	0.43
1:A:363:LYS:CE	1:A:363:LYS:HA	2.46	0.43
1:A:67:ILE:C	1:A:67:ILE:HD12	2.38	0.43
3:B:1436:P5S:H46A	3:B:1436:P5S:H50	1.74	0.43
1:A:104:LYS:HB3	1:A:104:LYS:HE2	1.70	0.43
3:A:1436:P5S:H20A	3:A:1436:P5S:H44A	2.01	0.42
1:B:232:MET:HE1	3:B:1436:P5S:H51A	2.01	0.42
1:B:225:ARG:O	1:B:232:MET:HA	2.19	0.42
1:A:378:ARG:HG3	1:A:378:ARG:HH11	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:224:VAL:CG2	1:B:232:MET:HE2	2.50	0.42
3:A:1436:P5S:H50	3:A:1436:P5S:H46A	1.86	0.41
1:B:96:ALA:O	1:B:105:ARG:HG3	2.19	0.41
1:B:67:ILE:O	1:B:67:ILE:HD12	2.20	0.41
1:A:227:ILE:CG2	1:A:232:MET:HE2	2.41	0.41
1:B:224:VAL:HG21	1:B:232:MET:HE2	2.01	0.41
1:B:67:ILE:HD12	1:B:67:ILE:C	2.42	0.41
1:A:257:GLY:HA3	1:A:260:PHE:H	1.85	0.40
1:B:92:PHE:CZ	1:B:111:LYS:HE2	2.56	0.40
1:A:96:ALA:O	1:A:105:ARG:HG3	2.22	0.40
1:B:186:LEU:CD2	5:B:1437:DTT:S1	3.10	0.40
1:B:72:PHE:HA	1:B:344:HIS:CE1	2.57	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)
6:B:2011:HOH:O	6:B:2195:HOH:O[4_556]	2.16	0.04

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	397/448 (89%)	377 (95%)	19 (5%)	1 (0%)	46	35
1	B	398/448 (89%)	380 (96%)	17 (4%)	1 (0%)	46	35
All	All	795/896 (89%)	757 (95%)	36 (4%)	2 (0%)	46	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	257	GLY

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Mol	Chain	Res	Type
1	B	261	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	360/405 (89%)	350 (97%)	10 (3%)	51	39
1	B	360/405 (89%)	350 (97%)	10 (3%)	51	39
All	All	720/810 (89%)	700 (97%)	20 (3%)	51	39

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	ASP
1	A	53	ASN
1	A	69	LEU
1	A	103	LEU
1	A	154	GLN
1	A	203	LEU
1	A	258	TYR
1	A	294	LEU
1	A	363	LYS
1	A	376	LEU
1	B	41	ASP
1	B	53	ASN
1	B	69	LEU
1	B	154	GLN
1	B	203	LEU
1	B	258	TYR
1	B	259	VAL
1	B	294	LEU
1	B	363	LYS
1	B	376	LEU

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

Mol	Chain	Res	Type
1	A	86	GLN
1	A	88	GLN
1	A	154	GLN
1	A	158	HIS
1	A	308	HIS
1	B	86	GLN
1	B	88	GLN
1	B	154	GLN
1	B	308	HIS
1	B	409	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	1435	-	4,4,4	0.20	0	6,6,6	0.16	0
3	P5S	A	1436	-	49,52,53	1.73	9 (18%)	49,59,60	1.49	7 (14%)
4	DTU	A	1437	-	7,7,7	1.39	1 (14%)	4,8,8	2.13	2 (50%)
2	SO4	B	1435	-	4,4,4	0.16	0	6,6,6	0.29	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	P5S	B	1436	-	50,53,53	1.57	7 (14%)	50,60,60	1.65	6 (12%)
5	DTT	B	1437	-	7,7,7	1.37	0	4,8,8	2.18	2 (50%)

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	SO4	A	1435	-	-	0/0/0/0	0/0/0/0
3	P5S	A	1436	-	-	0/54/58/59	0/0/0/0
4	DTU	A	1437	-	-	0/8/8/8	0/0/0/0
2	SO4	B	1435	-	-	0/0/0/0	0/0/0/0
3	P5S	B	1436	-	-	0/55/59/59	0/0/0/0
5	DTT	B	1437	-	-	0/8/8/8	0/0/0/0

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1436	P5S	P12-O15	-4.56	1.35	1.54
3	B	1436	P5S	P12-O15	-4.52	1.35	1.54
3	A	1436	P5S	C48-C46	-2.98	1.34	1.51
3	B	1436	P5S	C48-C46	-2.95	1.34	1.51
4	A	1437	DTU	O3-C3	-2.09	1.38	1.43
3	A	1436	P5S	P12-OG	-2.03	1.49	1.59
3	B	1436	P5S	O37-C2	2.15	1.52	1.46
3	A	1436	P5S	C20-C17	2.22	1.57	1.50
3	A	1436	P5S	O37-C2	2.23	1.52	1.46
3	B	1436	P5S	C1-C2	2.45	1.57	1.50
3	B	1436	P5S	O19-C17	3.13	1.42	1.33
3	A	1436	P5S	C1-C2	3.31	1.60	1.50
3	B	1436	P5S	C3-C2	3.35	1.60	1.50
3	A	1436	P5S	O19-C17	3.41	1.43	1.33
3	A	1436	P5S	C3-C2	3.73	1.61	1.50
3	B	1436	P5S	O37-C38	5.08	1.49	1.34
3	A	1436	P5S	O37-C38	5.09	1.49	1.34

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1436	P5S	O19-C1-C2	-4.10	97.65	108.69
3	B	1436	P5S	C22-C21-C20	-3.79	99.39	113.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1436	P5S	O37-C38-O47	-3.54	114.18	123.67
3	A	1436	P5S	O37-C38-O47	-3.22	115.03	123.67
3	A	1436	P5S	O19-C17-O18	-2.68	116.56	123.49
3	A	1436	P5S	OG-P12-O13	-2.37	100.41	109.62
3	B	1436	P5S	C25-C24-C23	-2.29	102.72	114.53
3	A	1436	P5S	C3-C2-C1	2.06	116.88	112.07
4	A	1437	DTU	O3-C3-C2	2.31	114.52	109.79
5	B	1437	DTT	O2-C2-C3	2.42	114.73	109.79
5	B	1437	DTT	C3-C4-S4	2.44	117.95	113.91
3	A	1436	P5S	C2-O37-C38	2.44	123.75	117.89
3	B	1436	P5S	C2-O37-C38	2.81	124.64	117.89
3	A	1436	P5S	O19-C17-C20	3.04	121.16	111.90
4	A	1437	DTU	C2-C1-S1	3.11	119.07	113.91
3	A	1436	P5S	O37-C38-C39	5.98	124.52	111.53
3	B	1436	P5S	O37-C38-C39	6.64	125.95	111.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1436	P5S	8	0
4	A	1437	DTU	3	0
3	B	1436	P5S	12	0
5	B	1437	DTT	5	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	399/448 (89%)	0.30	39 (9%)	10 15	17, 30, 62, 222	3 (0%)
1	B	400/448 (89%)	0.22	26 (6%)	22 32	15, 28, 60, 222	1 (0%)
All	All	799/896 (89%)	0.26	65 (8%)	15 23	15, 29, 61, 222	4 (0%)

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	258	TYR	10.8
1	A	258	TYR	10.8
1	A	259	VAL	9.9
1	B	259	VAL	8.7
1	A	52	LEU	8.4
1	B	366	ILE	7.3
1	A	257	GLY	6.8
1	B	260	PHE	6.7
1	B	208	GLY	6.0
1	A	51	ILE	5.9
1	B	369	GLY	5.6
1	A	298	ARG	5.3
1	A	48	HIS	5.1
1	A	230	GLY	5.1
1	A	55	ILE	5.0
1	A	56	SER	4.6
1	A	66	ARG	4.5
1	A	299	SER	4.3
1	A	296	GLN	4.3
1	A	431	ILE	4.3
1	A	54	ILE	3.6
1	A	359	ARG	3.5
1	A	260	PHE	3.4
1	A	300	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	261	GLY	3.4
1	B	257	GLY	3.2
1	B	207	ASP	3.0
1	A	262	THR	3.0
1	B	56	SER	3.0
1	A	143	PRO	3.0
1	A	261	GLY	2.9
1	A	58	LEU	2.8
1	B	180	ILE	2.8
1	A	370	VAL	2.8
1	B	35	ILE	2.7
1	B	370	VAL	2.7
1	A	309	LYS	2.6
1	A	433	SER	2.6
1	A	363	LYS	2.6
1	A	228	LEU	2.5
1	A	67	ILE	2.5
1	A	49	ASN	2.5
1	A	366	ILE	2.5
1	B	192	ALA	2.5
1	B	432	LYS	2.4
1	B	194	MET	2.4
1	A	369	GLY	2.4
1	B	397	ASP	2.3
1	B	229	PHE	2.3
1	B	48	HIS	2.3
1	A	341	VAL	2.3
1	A	229	PHE	2.3
1	A	297	PRO	2.2
1	A	45	GLU	2.2
1	B	237	GLY	2.2
1	B	179	VAL	2.2
1	A	57	GLN	2.2
1	B	45	GLU	2.2
1	B	66	ARG	2.1
1	A	302	LYS	2.1
1	B	368	ASP	2.1
1	B	209	ASN	2.1
1	A	217	LEU	2.0
1	B	196	GLY	2.0
1	A	237	GLY	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	SO4	A	1435	5/5	0.90	0.23	6.18	32,36,78,84	0
4	DTU	A	1437	8/8	0.85	0.30	5.29	48,68,84,96	0
5	DTT	B	1437	8/8	0.89	0.28	5.05	42,55,177,242	0
3	P5S	B	1436	54/54	0.96	0.17	0.79	13,21,36,74	0
2	SO4	B	1435	5/5	0.95	0.17	0.66	42,76,85,94	0
3	P5S	A	1436	53/54	0.93	0.14	0.28	16,28,37,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.