Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 02:09 AM GMT

PDB ID	:	2FVC
Title	:	Crystal structure of NS5B BK strain (delta 24) in complex with a 3-(1,1-Dioxo-2H-(1,2,4)-benzothiadiazin-3-yl)-4-hydroxy-2(1H)-quinolinone
Authors	:	Concha, N.O.; Wonacott, A.; Singh, O.
Deposited on	:	2006-01-30
Resolution	:	2.00 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

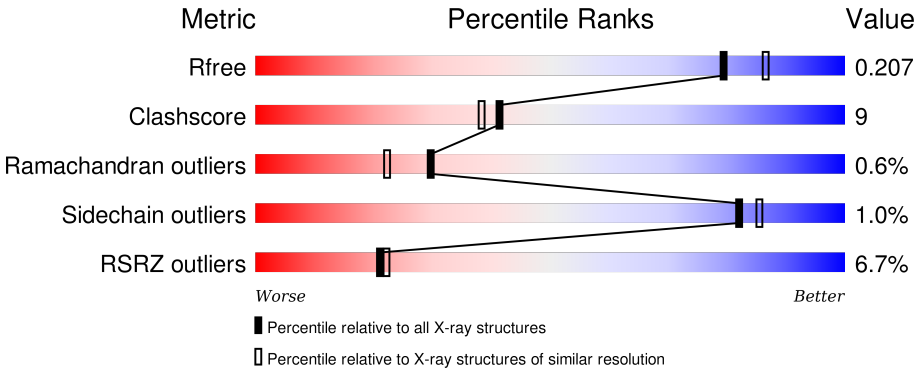
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 <=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	563	8% 83% 16%
1	B	563	6% 83% 16%

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	888	A	901	-	-	-	X
2	888	B	902	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9456 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.

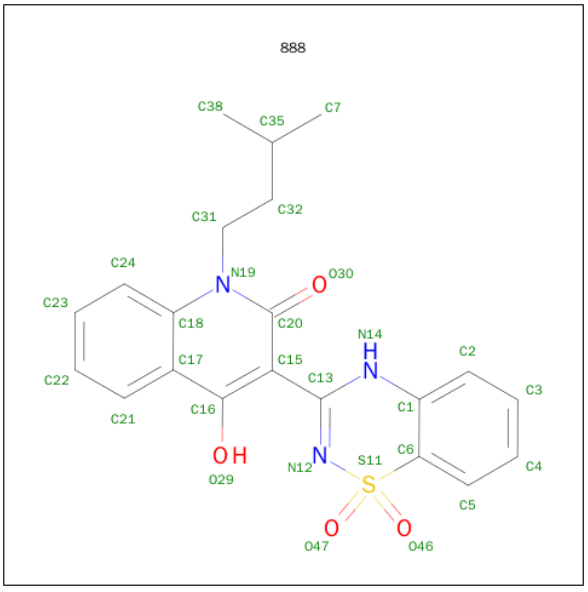
- Molecule 1 is a protein called polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	563	Total	C	N	O	S	14	2	0
			4376	2759	771	813	33			
1	B	563	Total	C	N	O	S	8	4	0
			4381	2764	771	813	33			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	329	VAL	THR	SEE REMARK 999	GB 12831193
B	329	VAL	THR	SEE REMARK 999	GB 12831193

- Molecule 2 is 3-(1,1-DIOXIDO-4H-1,2,4-BENZOTHIADIAZIN-3-YL)-4-HYDROXY-1-(3-METHYLBUTYL)QUINOLIN-2(1H)-ONE (three-letter code: 888) (formula: $C_{21}H_{21}N_3O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			29	21	3	4	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			29	21	3	4	1		

- Molecule 3 is water.

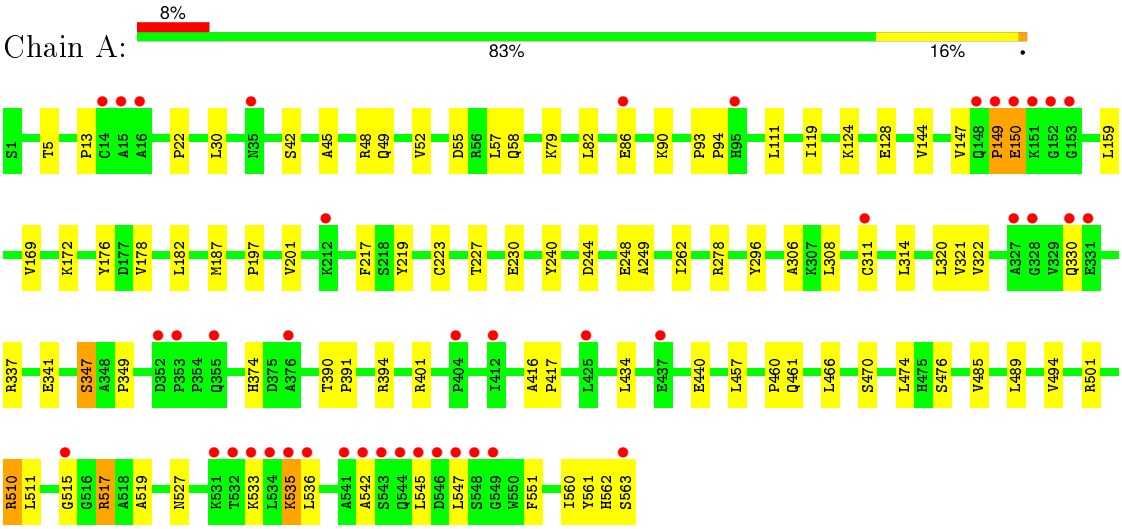
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	305	Total	O	0	0
			305	305		
3	B	336	Total	O	0	0
			336	336		

3 Residue-property plots (i)

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.

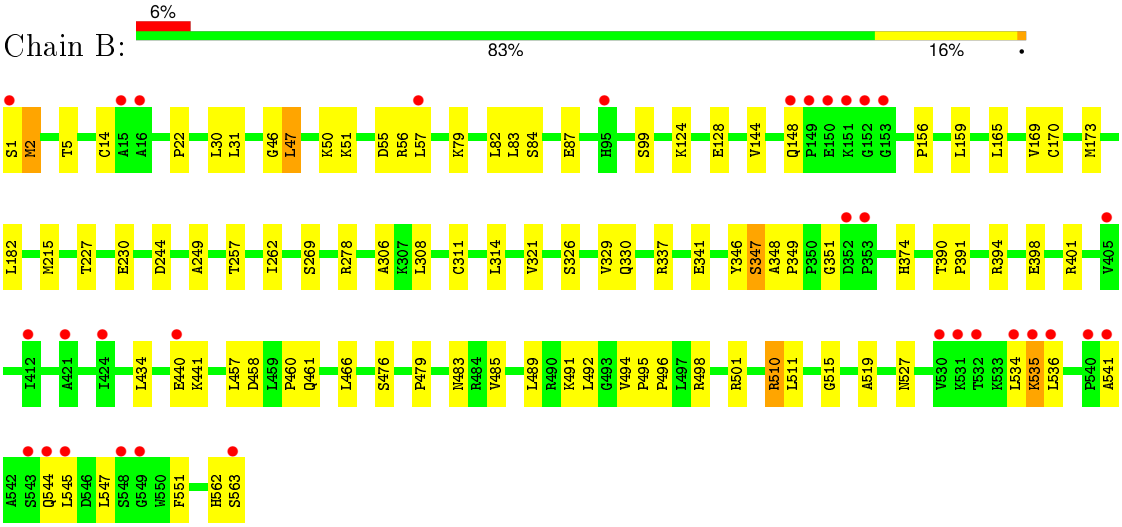
- Molecule 1: polyprotein

Chain A:



- Molecule 1: polyprotein

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	85.50Å 105.20Å 126.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 26.88 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.5 (20.00-2.00) 97.5 (26.88-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.41 (at 1.99Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.206 , 0.243 0.208 , 0.207	Depositor DCC
R_{free} test set	3743 reflections (4.95%)	DCC
Wilson B-factor (Å^2)	20.6	Xtriage
Anisotropy	0.437	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 53.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	2 of 75668 reflections (0.003%)	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	9456	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.88 % of the origin peak, indicating pseudo translational symmetry.The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.2407e-04.The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: 888

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	0.30	0/4482	0.57	0/6081
1	B	0.31	0/4497	0.59	2/6102 (0.0%)
All	All	0.30	0/8979	0.58	2/12183 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	351	GLY	N-CA-C	-5.57	99.17	113.10
1	B	31	LEU	CA-CB-CG	5.02	126.85	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4376	0	4396	75	0
1	B	4381	0	4408	75	0
2	A	29	0	21	0	0
2	B	29	0	21	0	0
3	A	305	0	0	7	0
3	B	336	0	0	4	0
All	All	9456	0	8846	150	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 9.

All (150) 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:5:THR:HG23	1:B:278:ARG:HH12	1.17	1.07
1:A:5:THR:HG23	1:A:278:ARG:HH12	1.24	0.99
1:A:535:LYS:H	1:A:535:LYS:HD3	1.48	0.78
1:A:330:GLN:HB2	3:A:1201:HOH:O	1.88	0.72
1:B:308:LEU:HB3	1:B:311:CYS:SG	2.29	0.72
1:A:308:LEU:HB3	1:A:311:CYS:SG	2.30	0.71
1:A:5:THR:HG21	1:A:278:ARG:HH22	1.55	0.71
1:A:347:SER:O	1:A:349:PRO:HD3	1.91	0.70
1:B:510:ARG:HH11	1:B:510:ARG:HG2	1.58	0.68
1:B:5:THR:HG21	1:B:278:ARG:HH22	1.57	0.67
1:B:5:THR:HG23	1:B:278:ARG:NH1	2.00	0.67
1:A:149:PRO:HG2	1:A:150:GLU:H	1.58	0.67
1:B:337:ARG:O	1:B:341:GLU:HG3	1.95	0.66
1:A:517:ARG:HG2	1:A:517:ARG:HH11	1.60	0.66
1:B:527:ASN:HD21	1:B:534:LEU:H	1.45	0.65
1:B:535:LYS:HG3	1:B:536:LEU:H	1.61	0.65
1:B:1:SER:HB3	3:B:1214:HOH:O	1.96	0.65
1:A:178:VAL:HG23	3:A:1017:HOH:O	1.95	0.65
1:B:440:GLU:HG2	1:B:457:LEU:HD12	1.78	0.64
1:A:13:PRO:HG3	1:A:42:SER:OG	1.98	0.64
1:A:440:GLU:HG2	1:A:457:LEU:HD12	1.80	0.63
1:B:510:ARG:HG2	3:B:1205:HOH:O	1.99	0.63
1:B:182:LEU:HD23	1:B:182:LEU:C	2.19	0.63
1:A:466:LEU:HD22	1:A:551:PHE:HE2	1.62	0.62
1:A:176:TYR:OH	1:A:562:HIS:HE1	1.82	0.61
1:A:124:LYS:O	1:A:128:GLU:HG3	2.02	0.59
1:B:346:TYR:O	1:B:347:SER:HB3	2.01	0.59
1:A:22:PRO:HG2	1:A:401:ARG:HG3	1.85	0.59
1:A:52:VAL:HG12	1:A:223[B]:CYS:SG	2.43	0.58
1:B:170:CYS:HA	1:B:173:MET:HE3	1.85	0.58
1:A:337:ARG:O	1:A:341:GLU:HG3	2.03	0.57
1:A:182:LEU:HD23	1:A:182:LEU:C	2.24	0.57
1:A:510:ARG:HH11	1:A:510:ARG:HG2	1.70	0.56
1:B:169:VAL:HG12	1:B:173:MET:HE2	1.87	0.56
1:A:347:SER:C	1:A:349:PRO:HD3	2.25	0.56
1:A:306:ALA:HB3	1:A:308:LEU:HD13	1.88	0.56
1:B:50:LYS:HD2	1:B:50:LYS:N	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:LEU:HD22	1:B:51:LYS:NZ	2.20	0.56
1:B:314:LEU:HB3	1:B:321:VAL:HG12	1.89	0.55
1:A:308:LEU:CB	1:A:311:CYS:SG	2.94	0.55
1:B:124:LYS:O	1:B:128:GLU:HG3	2.06	0.55
1:B:515:GLY:HA2	1:B:519:ALA:HB2	1.89	0.55
1:B:510:ARG:NH1	1:B:510:ARG:HG2	2.21	0.55
1:A:314:LEU:HB3	1:A:321:VAL:CG1	2.36	0.55
1:A:314:LEU:HB3	1:A:321:VAL:HG12	1.89	0.54
1:A:501:ARG:NH1	1:A:501:ARG:HB2	2.22	0.54
1:B:227:THR:HB	1:B:347:SER:O	2.07	0.54
1:A:517:ARG:HG2	1:A:517:ARG:NH1	2.23	0.54
1:A:230:GLU:HG3	1:A:262:ILE:HG23	1.90	0.54
1:B:348:ALA:N	1:B:349:PRO:HD3	2.23	0.54
1:A:501:ARG:HB2	1:A:501:ARG:HH11	1.73	0.53
1:A:5:THR:HG23	1:A:278:ARG:NH1	2.08	0.53
1:A:460:PRO:HB2	1:A:461:GLN:HE21	1.74	0.53
1:A:535:LYS:HE3	3:A:1204:HOH:O	2.09	0.53
1:A:390:THR:HB	1:A:391:PRO:HD3	1.91	0.53
1:B:374:HIS:HD2	1:B:476:SER:HB2	1.74	0.53
1:A:434:LEU:HD21	1:A:511:LEU:HD23	1.90	0.53
1:A:240:TYR:OH	3:A:1193:HOH:O	2.14	0.53
1:B:82:LEU:HD13	1:B:249:ALA:HB2	1.89	0.52
1:B:390:THR:HB	1:B:391:PRO:HD3	1.91	0.52
1:B:541:ALA:O	1:B:544:GLN:HG2	2.10	0.52
1:A:230:GLU:HG3	1:A:262:ILE:CG2	2.40	0.52
1:B:215:MET:HB2	1:B:326:SER:HB2	1.91	0.52
1:B:230:GLU:HG3	1:B:262[A]:ILE:CG2	2.39	0.52
1:A:248:GLU:HB3	3:A:1075:HOH:O	2.09	0.52
1:A:535:LYS:HG2	3:A:1204:HOH:O	2.09	0.51
1:B:308:LEU:CB	1:B:311:CYS:SG	2.97	0.51
1:B:47:LEU:HD13	1:B:156:PRO:HG3	1.92	0.51
1:B:394:ARG:O	1:B:398:GLU:HG3	2.10	0.51
1:A:5:THR:CG2	1:A:278:ARG:HH12	2.11	0.51
1:A:144:VAL:HB	1:A:394:ARG:HG2	1.92	0.51
1:B:314:LEU:HB3	1:B:321:VAL:CG1	2.42	0.50
1:B:22:PRO:HG2	1:B:401:ARG:HG3	1.92	0.50
1:B:535:LYS:O	1:B:536:LEU:HB2	2.12	0.50
1:B:466:LEU:HD22	1:B:551:PHE:HE2	1.75	0.50
1:A:545:LEU:HB3	1:A:547:LEU:CD1	2.42	0.50
1:B:306:ALA:HB3	1:B:308:LEU:HD13	1.94	0.49
1:A:461:GLN:HB2	1:A:545:LEU:HD11	1.92	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:230:GLU:HG3	1:B:262[A]:ILE:HG23	1.94	0.49
1:B:47:LEU:HD22	1:B:51:LYS:HZ3	1.78	0.49
1:B:5:THR:CG2	1:B:278:ARG:HH22	2.26	0.49
1:A:545:LEU:HB3	1:A:547:LEU:HD13	1.94	0.49
1:B:329:VAL:HG23	1:B:330:GLN:N	2.28	0.49
1:B:545:LEU:HB3	1:B:547:LEU:CD1	2.43	0.49
1:B:144:VAL:HB	1:B:394:ARG:HG2	1.94	0.49
1:A:57:LEU:HD12	1:A:57:LEU:N	2.28	0.48
1:B:170:CYS:HA	1:B:173:MET:CE	2.44	0.48
1:B:2:MET:H	1:B:55:ASP:HA	1.77	0.48
1:B:169:VAL:HG12	1:B:173:MET:CE	2.43	0.48
1:A:510:ARG:NH1	1:A:510:ARG:HG2	2.28	0.48
1:A:86:GLU:HG3	1:A:111:LEU:HD11	1.94	0.48
1:A:48:ARG:HG2	1:A:159:LEU:HD13	1.96	0.48
1:A:440:GLU:HG2	1:A:457:LEU:CD1	2.43	0.48
1:B:440:GLU:HG2	1:B:457:LEU:CD1	2.44	0.47
1:B:182:LEU:HD23	1:B:182:LEU:O	2.14	0.47
1:A:172:LYS:HE3	1:A:560:ILE:HD13	1.97	0.47
1:B:491:LYS:HE3	1:B:492:LEU:CD1	2.45	0.47
1:A:119:ILE:HD13	1:A:169:VAL:HG11	1.95	0.47
1:A:374:HIS:HD2	1:A:476:SER:HB2	1.79	0.47
1:A:176:TYR:OH	1:A:562:HIS:CE1	2.64	0.46
1:B:434:LEU:HD21	1:B:511:LEU:HD23	1.96	0.46
1:B:55:ASP:OD2	1:B:57:LEU:HD11	2.16	0.46
1:A:86:GLU:O	1:A:90:LYS:HG2	2.16	0.46
1:B:56:ARG:C	1:B:57:LEU:HD12	2.36	0.46
1:A:561:TYR:CE2	1:A:563:SER:HB2	2.51	0.46
1:A:535:LYS:CD	1:A:535:LYS:H	2.24	0.46
1:A:485:VAL:O	1:A:489:LEU:HG	2.16	0.45
1:B:491:LYS:HE3	1:B:492:LEU:HD11	1.98	0.45
1:B:461:GLN:HB2	1:B:545:LEU:HD11	1.98	0.45
1:A:82:LEU:HD13	1:A:249:ALA:HB2	1.99	0.45
1:A:45:ALA:O	1:A:49:GLN:HG3	2.17	0.45
1:B:14[A]:CYS:SG	1:B:269:SER:HB2	2.57	0.44
1:A:501:ARG:NH1	3:A:1195:HOH:O	2.50	0.44
1:A:227:THR:HB	1:A:347:SER:O	2.16	0.44
1:B:84:SER:OG	1:B:87:GLU:HG3	2.17	0.44
1:B:1:SER:N	3:B:1226:HOH:O	2.35	0.44
1:A:219:TYR:HB3	1:A:320:LEU:HD23	1.98	0.44
1:A:515:GLY:HA2	1:A:519:ALA:HB2	2.00	0.43
1:B:99:SER:HB2	1:B:165:LEU:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:LYS:HA	1:A:244:ASP:HB3	1.99	0.43
1:B:479:PRO:O	1:B:483:ASN:ND2	2.52	0.43
1:B:441:LYS:HD3	3:B:1148:HOH:O	2.17	0.43
1:B:46:GLY:O	1:B:50:LYS:HD3	2.19	0.43
1:A:5:THR:CG2	1:A:278:ARG:HH22	2.29	0.42
1:A:416:ALA:N	1:A:417:PRO:CD	2.82	0.42
1:B:257:THR:O	1:B:262[B]:ILE:HG23	2.19	0.42
1:B:562:HIS:O	1:B:563:SER:C	2.57	0.42
1:B:498:ARG:HA	1:B:501:ARG:NH1	2.34	0.42
1:A:561:TYR:CZ	1:A:563:SER:HB2	2.54	0.42
1:A:93:PRO:HA	1:A:94:PRO:HD3	1.98	0.42
1:A:470:SER:O	1:A:474:LEU:HG	2.19	0.42
1:A:58:GLN:HG2	1:A:347:SER:HB3	2.01	0.42
1:B:458:ASP:HA	1:B:461:GLN:NE2	2.35	0.42
1:A:197:PRO:O	1:A:201:VAL:HG23	2.20	0.42
1:A:461:GLN:HB3	1:A:542:ALA:HA	2.01	0.42
1:B:461:GLN:N	1:B:461:GLN:OE1	2.48	0.41
1:A:187:MET:HE3	1:A:296:TYR:HB2	2.01	0.41
1:A:217:PHE:CE1	1:A:322:VAL:HB	2.55	0.41
1:B:79:LYS:HA	1:B:244:ASP:HB3	2.02	0.41
1:A:30:LEU:O	1:A:494:VAL:HG22	2.21	0.41
1:B:485:VAL:O	1:B:489:LEU:HG	2.20	0.41
1:B:83:LEU:HB2	1:B:173:MET:HA	2.01	0.41
1:B:169:VAL:O	1:B:173:MET:HE3	2.20	0.41
1:A:55:ASP:OD2	1:A:57:LEU:HD11	2.21	0.41
1:B:515:GLY:CA	1:B:519:ALA:HB2	2.51	0.41
1:B:460:PRO:HB2	1:B:461:GLN:OE1	2.21	0.41
1:B:30:LEU:O	1:B:494:VAL:HG22	2.21	0.41
1:A:149:PRO:HG2	1:A:150:GLU:N	2.32	0.40
1:B:374:HIS:CD2	1:B:476:SER:HB2	2.56	0.40
1:B:495:PRO:HA	1:B:496:PRO:HD3	1.94	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	563/563 (100%)	542 (96%)	17 (3%)	4 (1%)	26	19
1	B	565/563 (100%)	546 (97%)	16 (3%)	3 (0%)	34	26
All	All	1128/1126 (100%)	1088 (96%)	33 (3%)	7 (1%)	30	22

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	149	PRO
1	B	347	SER
1	B	535	LYS
1	A	533	LYS
1	A	536	LEU
1	B	2	MET
1	A	147	VAL

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	481/479 (100%)	475 (99%)	6 (1%)	78	81
1	B	483/479 (101%)	479 (99%)	4 (1%)	86	89
All	All	964/958 (101%)	954 (99%)	10 (1%)	82	85

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	150	GLU
1	A	347	SER
1	A	510	ARG
1	A	517	ARG
1	A	527	ASN
1	A	535	LYS

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Mol	Chain	Res	Type
1	B	47	LEU
1	B	148	GLN
1	B	159	LEU
1	B	510	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
1	A	49	GLN
1	A	273	ASN
1	A	309	GLN
1	A	374	HIS
1	A	461	GLN
1	A	544	GLN
1	A	562	HIS
1	B	49	GLN
1	B	120	HIS
1	B	273	ASN
1	B	374	HIS
1	B	483	ASN
1	B	527	ASN
1	B	544	GLN

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

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	888	A	901	-	31,32,32	1.64	7 (22%)	36,48,48	2.11	7 (19%)
2	888	B	902	-	31,32,32	1.69	6 (19%)	36,48,48	2.01	7 (19%)

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	888	A	901	-	-	0/8/24/24	0/4/4/4
2	888	B	902	-	-	0/8/24/24	0/4/4/4

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	902	888	C1-N14	-2.94	1.34	1.39
2	A	901	888	C1-N14	-2.55	1.35	1.39
2	A	901	888	C5-C6	2.01	1.41	1.39
2	B	902	888	C16-C17	2.13	1.48	1.43
2	A	901	888	C16-C17	2.14	1.48	1.43
2	B	902	888	C23-C24	2.24	1.41	1.36
2	A	901	888	C18-N19	2.28	1.43	1.40
2	A	901	888	C23-C24	2.32	1.42	1.36
2	B	902	888	C18-N19	2.48	1.43	1.40
2	A	901	888	C6-S11	3.24	1.78	1.74
2	B	902	888	C6-S11	3.33	1.78	1.74
2	A	901	888	C20-C15	4.09	1.53	1.44
2	B	902	888	C20-C15	4.15	1.53	1.44

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	902	888	O46-S11-N12	-2.50	105.86	109.15
2	A	901	888	O46-S11-N12	-2.48	105.88	109.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	902	888	C15-C13-N14	-2.07	115.53	118.56
2	A	901	888	O47-S11-O46	2.00	118.26	115.75
2	A	901	888	C31-N19-C18	2.00	121.01	119.06
2	B	902	888	C21-C17-C18	2.06	120.46	118.17
2	A	901	888	C21-C17-C18	2.07	120.47	118.17
2	A	901	888	C1-N14-C13	2.23	129.51	123.92
2	B	902	888	O47-S11-O46	2.28	118.60	115.75
2	B	902	888	C1-N14-C13	2.39	129.90	123.92
2	A	901	888	C15-C13-N12	3.49	122.79	118.74
2	B	902	888	C15-C13-N12	4.01	123.39	118.74
2	B	902	888	C32-C31-N19	8.64	119.59	112.29
2	A	901	888	C32-C31-N19	9.89	120.65	112.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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^{th} 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	563/563 (100%)	0.40	43 (7%)	17	18	10, 21, 42, 61	3 (0%)
1	B	563/563 (100%)	0.32	32 (5%)	27	29	10, 19, 37, 60	2 (0%)
All	All	1126/1126 (100%)	0.36	75 (6%)	21	22	10, 20, 40, 61	5 (0%)

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	149	PRO	7.2
1	A	150	GLU	7.2
1	B	532	THR	7.2
1	B	152	GLY	7.2
1	B	149	PRO	6.9
1	A	152	GLY	6.9
1	A	534	LEU	6.6
1	B	150	GLU	6.2
1	B	151	LYS	6.1
1	A	151	LYS	6.0
1	B	563	SER	5.7
1	B	549	GLY	5.1
1	B	544	GLN	4.7
1	B	534	LEU	4.4
1	A	548	SER	4.2
1	A	541	ALA	4.2
1	B	148	GLN	4.2
1	A	549	GLY	4.1
1	B	531	LYS	4.1
1	B	153	GLY	4.1
1	B	548	SER	4.0
1	A	531	LYS	3.8
1	B	1	SER	3.6
1	A	544	GLN	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	16	ALA	3.5
1	A	148	GLN	3.4
1	B	535	LYS	3.4
1	B	541	ALA	3.4
1	A	412	ILE	3.4
1	A	546	ASP	3.3
1	B	543	SER	3.3
1	A	532	THR	3.3
1	A	542	ALA	3.1
1	A	331	GLU	3.0
1	B	16	ALA	3.0
1	A	535	LYS	3.0
1	B	530	VAL	3.0
1	B	412	ILE	3.0
1	B	57	LEU	3.0
1	B	545	LEU	2.9
1	B	540	PRO	2.9
1	A	536	LEU	2.9
1	A	563	SER	2.9
1	A	327	ALA	2.8
1	A	545	LEU	2.8
1	A	353	PRO	2.8
1	A	355	GLN	2.8
1	A	330	GLN	2.7
1	A	515	GLY	2.7
1	A	212	LYS	2.7
1	B	15	ALA	2.7
1	A	404	PRO	2.7
1	B	352	ASP	2.7
1	A	311	CYS	2.6
1	A	547	LEU	2.6
1	A	376	ALA	2.6
1	B	95	HIS	2.5
1	A	153	GLY	2.4
1	A	352	ASP	2.4
1	A	328	GLY	2.4
1	A	35	ASN	2.3
1	A	14	CYS	2.3
1	A	425	LEU	2.2
1	B	405	VAL	2.2
1	A	437	GLU	2.2
1	B	421	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	95	HIS	2.1
1	A	15	ALA	2.1
1	A	543	SER	2.1
1	B	536	LEU	2.1
1	B	353	PRO	2.1
1	B	440	GLU	2.1
1	B	424	ILE	2.0
1	A	533	LYS	2.0
1	A	86	GLU	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^{th} 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^2)	Q<0.9
2	888	A	901	29/29	0.89	0.17	2.88	20,23,26,33	0
2	888	B	902	29/29	0.90	0.17	2.55	17,20,22,27	0

6.5 Other polymers (i)

There are no such residues in this entry.