Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 02:06 PM GMT

PDB ID	:	3W5E
Title	:	Crystal structure of phosphodiesterase 4B in complex with compound 31e
Authors	:	Takahashi, M.; Hanzawa, H.
Deposited on	:	2013-01-28
Resolution	:	2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
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:

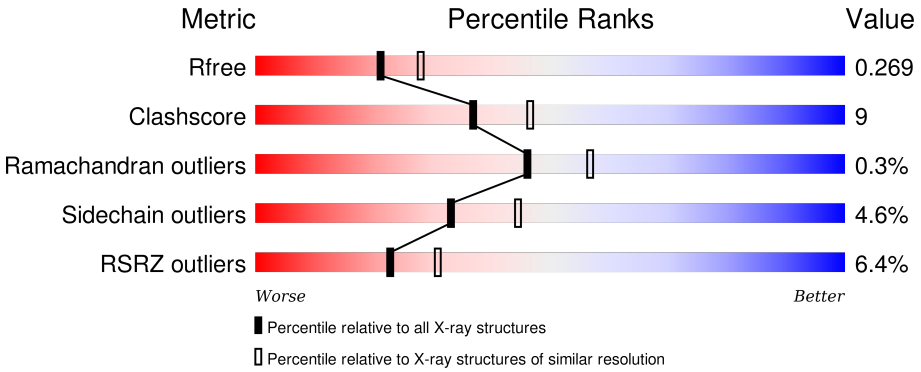
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	377	6% 75% 14% • • 8%
1	B	377	5% 68% 22% • 8%

2 Entry composition

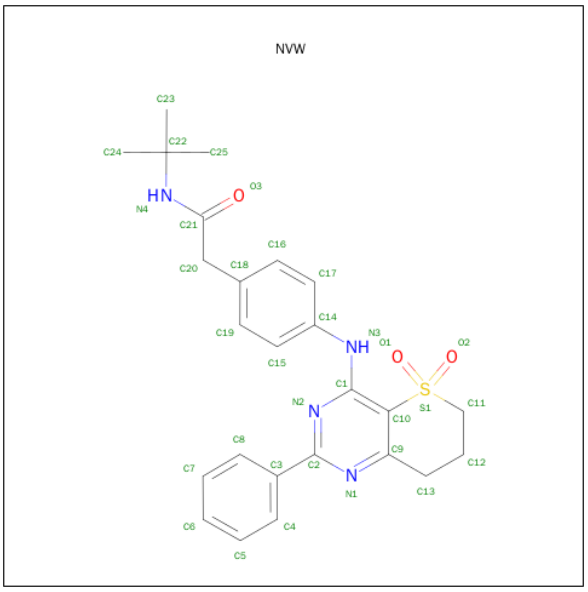
There are 5 unique types of molecules in this entry. The entry contains 5994 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 cAMP-specific 3',5'-cyclic phosphodiesterase 4B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	345	Total	C	N	O	S	0	0	0
			2792	1763	471	537	21			
1	B	345	Total	C	N	O	S	0	0	0
			2792	1763	471	537	21			

- Molecule 2 is N-TERT-BUTYL-2-{4-[(5,5-DIOXIDO-2-PHENYL-7,8-DIHYDRO-6H-THIOPYRANO[3,2-D]PYRIMIDIN-4-YL)AMINO]PHENYL}ACETAMIDE (three-letter code: NVW) (formula: $C_{25}H_{28}N_4O_3S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			33	25	4	3	1		
2	B	1	Total	C	N	O	S	0	0
			33	25	4	3	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Ca 1 1	0	0
4	A	1	Total Ca 1 1	0	0

- Molecule 5 is water.

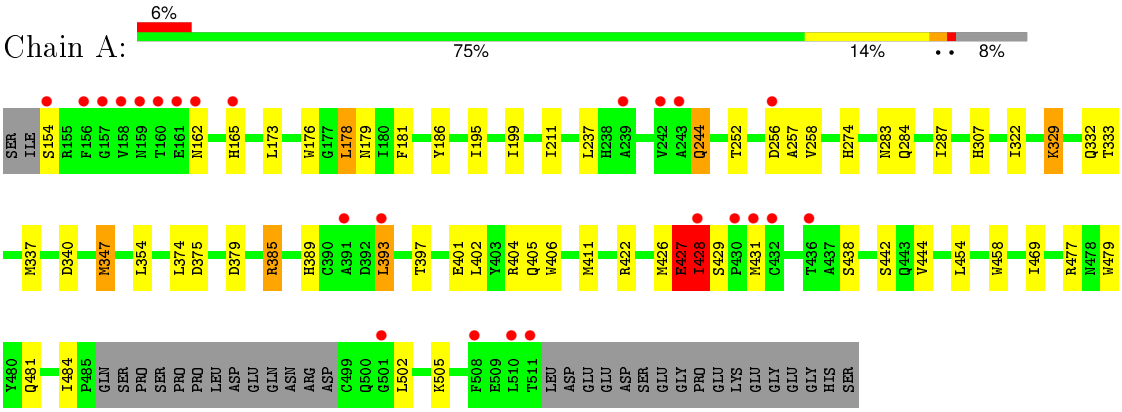
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	172	Total O 172 172	0	0
5	B	168	Total O 168 168	0	0

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.

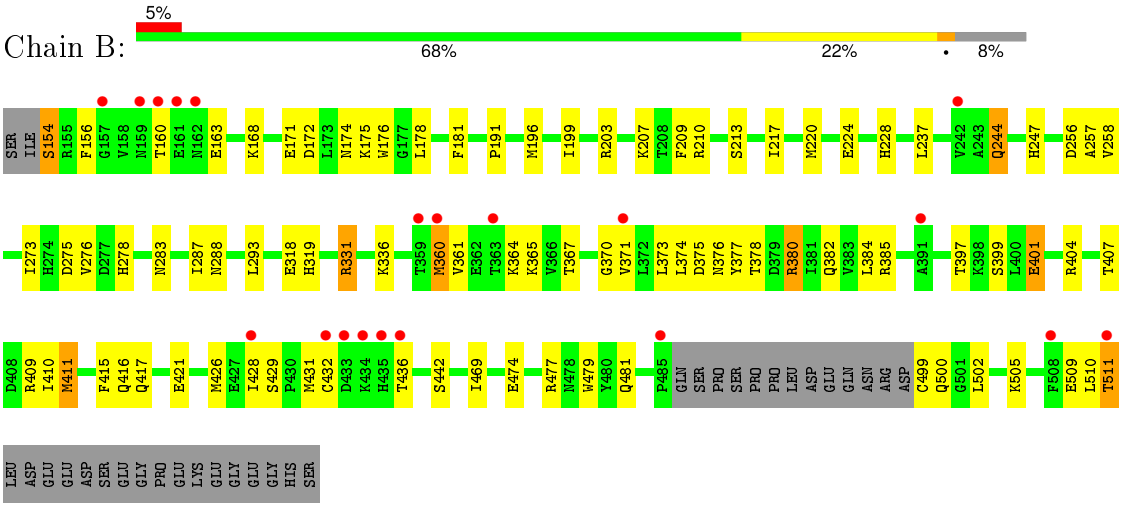
- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4B

Chain A:



- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4B

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	81.06Å 158.19Å 58.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.77 – 2.30 35.19 – 2.29	Depositor EDS
% Data completeness (in resolution range)	88.5 (19.77-2.30) 88.4 (35.19-2.29)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.10 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.182 , 0.269 0.182 , 0.269	Depositor DCC
R_{free} test set	3027 reflections (11.15%)	DCC
Wilson B-factor (Å^2)	35.5	Xtriage
Anisotropy	0.326	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 33.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Outliers	0 of 30497 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5994	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.48% of the height of the origin peak. No significant pseudotranslation is detected.*

[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: NVW, ZN, CA

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	0.69	4/2849 (0.1%)	0.71	0/3859
1	B	0.69	2/2849 (0.1%)	0.73	1/3859 (0.0%)
All	All	0.69	6/5698 (0.1%)	0.72	1/7718 (0.0%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	479	TRP	CD2-CE2	5.73	1.48	1.41
1	A	458	TRP	CD2-CE2	5.35	1.47	1.41
1	A	176	TRP	CD2-CE2	5.33	1.47	1.41
1	B	176	TRP	CD2-CE2	5.31	1.47	1.41
1	A	479	TRP	CD2-CE2	5.19	1.47	1.41
1	A	406	TRP	CD2-CE2	5.09	1.47	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	409	ARG	NE-CZ-NH2	-7.86	116.37	120.30

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	2792	0	2710	44	0
1	B	2792	0	2710	61	0
2	A	33	0	28	1	0
2	B	33	0	28	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	172	0	0	7	0
5	B	168	0	0	6	0
All	All	5994	0	5476	105	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 (105) 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:367:THR:HG22	1:B:371:VAL:H	1.20	0.99
1:A:428:ILE:O	1:A:428:ILE:HD13	1.63	0.96
1:A:427:GLU:O	1:A:427:GLU:HG2	1.65	0.94
1:A:284:GLN:OE1	5:A:1089:HOH:O	1.96	0.83
1:B:331:ARG:NH1	5:B:1072:HOH:O	2.13	0.81
1:B:203:ARG:HE	1:B:247:HIS:HD2	1.26	0.80
1:B:367:THR:HG22	1:B:371:VAL:N	1.98	0.78
1:B:367:THR:CG2	1:B:371:VAL:H	1.98	0.77
1:B:511:THR:CG2	5:B:1107:HOH:O	2.35	0.74
1:A:428:ILE:CD1	1:A:428:ILE:O	2.35	0.73
1:B:411:MET:CE	1:B:502:LEU:HD13	2.20	0.72
1:B:411:MET:HE1	1:B:502:LEU:HD13	1.72	0.70
1:A:442:SER:HB3	2:A:901:NVW:H3	1.72	0.70
1:A:427:GLU:O	1:A:427:GLU:CG	2.40	0.69
1:B:367:THR:HG23	1:B:370:GLY:H	1.59	0.67
1:B:511:THR:HG21	5:B:1107:HOH:O	1.95	0.67
1:B:431:MET:HE2	5:B:1085:HOH:O	1.93	0.67
1:B:178:LEU:HD22	1:B:244:GLN:HG3	1.76	0.67
1:A:256:ASP:O	1:A:257:ALA:HB3	1.94	0.67
1:B:203:ARG:NE	1:B:247:HIS:HD2	1.94	0.65
1:A:427:GLU:O	1:A:429:SER:N	2.28	0.65
1:B:196:MET:HG3	1:B:220:MET:HE1	1.80	0.63
1:A:283:ASN:O	1:A:287:ILE:HG13	1.98	0.63
1:B:283:ASN:O	1:B:287:ILE:HG13	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:LEU:HD22	1:A:244:GLN:HG3	1.82	0.62
1:A:252:THR:HG21	5:A:1165:HOH:O	1.99	0.61
1:B:160:THR:HA	1:B:163:GLU:HG3	1.83	0.60
1:B:509:GLU:O	1:B:510:LEU:C	2.38	0.59
1:B:196:MET:HG3	1:B:220:MET:CE	2.32	0.59
1:B:288:ASN:HD22	1:B:426:MET:HE1	1.68	0.59
1:B:442:SER:HB3	2:B:901:NVW:H3	1.84	0.58
1:B:288:ASN:HD22	1:B:426:MET:CE	2.16	0.57
1:A:428:ILE:CG1	1:A:428:ILE:O	2.47	0.57
1:A:431:MET:HE3	1:A:505:LYS:HD3	1.85	0.57
1:A:428:ILE:HG23	1:A:428:ILE:O	2.05	0.57
1:A:179:ASN:ND2	1:A:402:LEU:HD21	2.19	0.57
1:A:404:ARG:HD2	5:A:1064:HOH:O	2.05	0.56
1:A:426:MET:O	1:A:427:GLU:C	2.44	0.56
1:B:505:LYS:O	1:B:509:GLU:HG3	2.05	0.55
1:A:374:LEU:HD22	1:A:379:ASP:HB3	1.88	0.55
1:A:477:ARG:O	1:A:481:GLN:HB2	2.07	0.55
1:B:336:LYS:HE3	1:B:377:TYR:CZ	2.42	0.55
1:A:411:MET:HE2	1:A:502:LEU:HD11	1.91	0.53
1:B:213:SER:O	1:B:217:ILE:HG12	2.08	0.53
1:A:244:GLN:NE2	5:A:1170:HOH:O	2.38	0.52
1:B:415:PHE:HD1	1:B:431:MET:HE3	1.76	0.51
1:B:382:GLN:HG3	1:B:385:ARG:HH11	1.75	0.51
1:B:428:ILE:HG12	1:B:429:SER:N	2.24	0.51
1:B:360:MET:HG3	1:B:361:VAL:N	2.26	0.51
1:A:347:MET:CE	1:A:393:LEU:HD11	2.41	0.51
1:A:256:ASP:O	1:A:257:ALA:CB	2.59	0.50
1:B:417:GLN:O	1:B:421:GLU:HG3	2.11	0.50
1:A:438:SER:HB2	5:A:1085:HOH:O	2.10	0.50
1:B:411:MET:HE2	1:B:502:LEU:HD13	1.93	0.50
1:B:380:ARG:O	1:B:384:LEU:HG	2.13	0.49
1:B:376:ASN:HD22	1:B:378:THR:H	1.59	0.49
1:B:399:SER:HB3	1:B:401:GLU:OE2	2.12	0.48
1:A:181:PHE:CD1	1:A:237:LEU:HD21	2.49	0.48
1:B:365:LYS:HB2	1:B:373:LEU:HB2	1.95	0.48
1:A:165:HIS:HB3	1:A:186:TYR:CE2	2.49	0.47
1:B:168:LYS:HD3	1:B:171:GLU:OE1	2.14	0.47
1:B:336:LYS:HE3	1:B:377:TYR:OH	2.14	0.47
1:A:389:HIS:CE1	1:A:393:LEU:HD22	2.50	0.47
1:A:258:VAL:HG11	1:A:374:LEU:HD12	1.96	0.46
1:B:174:ASN:OD1	1:B:247:HIS:HE1	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:411:MET:HE2	1:B:502:LEU:CD1	2.46	0.46
1:B:477:ARG:HH12	1:B:481:GLN:NE2	2.13	0.45
1:B:181:PHE:CD1	1:B:237:LEU:HD21	2.52	0.44
1:B:228:HIS:HD2	1:B:278:HIS:O	2.00	0.44
1:B:365:LYS:HE2	1:B:365:LYS:HA	1.99	0.44
1:B:367:THR:CG2	1:B:370:GLY:H	2.27	0.44
1:A:211:ILE:HG12	1:A:322:ILE:O	2.17	0.44
1:A:411:MET:HE3	1:A:411:MET:HB3	1.87	0.44
1:B:209:PHE:O	1:B:210:ARG:C	2.55	0.44
1:A:426:MET:HG2	5:A:1156:HOH:O	2.18	0.44
1:B:401:GLU:H	1:B:401:GLU:HG3	1.39	0.44
1:A:195:ILE:O	1:A:199:ILE:HG13	2.18	0.44
1:B:172:ASP:HA	1:B:175:LYS:HD2	2.00	0.43
1:A:397:THR:HB	1:A:469:ILE:HG23	1.99	0.43
1:B:415:PHE:HA	1:B:431:MET:HE1	2.01	0.43
1:B:318:GLU:HG3	1:B:319:HIS:CE1	2.53	0.43
1:A:354:LEU:HD13	1:A:454:LEU:HA	2.01	0.43
1:A:347:MET:HE3	1:A:393:LEU:HD11	2.00	0.43
1:A:274:HIS:O	1:A:307:HIS:CD2	2.72	0.43
1:B:275:ASP:O	1:B:278:HIS:HB2	2.19	0.43
1:B:426:MET:HB3	1:B:426:MET:HE2	1.57	0.42
1:A:244:GLN:HE21	1:A:244:GLN:HB3	1.64	0.42
1:B:397:THR:HB	1:B:469:ILE:HG23	2.00	0.42
1:B:364:LYS:NZ	5:B:1036:HOH:O	2.53	0.42
1:B:191:PRO:HD2	1:B:224:GLU:OE2	2.18	0.42
1:A:401:GLU:O	1:A:405:GLN:HG3	2.19	0.42
1:A:444:VAL:HG21	1:A:484:ILE:HD11	2.02	0.41
1:A:422:ARG:NH2	5:A:1027:HOH:O	2.34	0.41
1:A:340:ASP:OD2	1:A:385:ARG:NE	2.48	0.41
1:B:273:ILE:O	1:B:276:VAL:HG12	2.20	0.41
1:B:154:SER:HB3	1:B:156:PHE:H	1.84	0.41
1:A:329:LYS:HA	1:A:332:GLN:HG2	2.03	0.41
1:B:407:THR:O	1:B:410:ILE:HG22	2.21	0.41
1:B:199:ILE:HD11	1:B:244:GLN:HA	2.02	0.41
1:B:331:ARG:NH2	5:B:1069:HOH:O	2.47	0.41
1:B:258:VAL:HG11	1:B:374:LEU:HD12	2.03	0.41
1:A:333:THR:O	1:A:337:MET:HG3	2.21	0.41
1:A:173:LEU:HA	1:A:178:LEU:HD13	2.03	0.40
1:B:256:ASP:O	1:B:257:ALA:HB3	2.22	0.40
1:B:510:LEU:HD12	1:B:510:LEU:HA	1.93	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	341/377 (90%)	328 (96%)	11 (3%)	2 (1%)	30 36
1	B	341/377 (90%)	329 (96%)	12 (4%)	0	100 100
All	All	682/754 (90%)	657 (96%)	23 (3%)	2 (0%)	46 57

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	427	GLU
1	A	428	ILE

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	313/342 (92%)	302 (96%)	11 (4%)	43 58
1	B	313/342 (92%)	295 (94%)	18 (6%)	25 33
All	All	626/684 (92%)	597 (95%)	29 (5%)	33 44

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	154	SER
1	A	162	ASN
1	A	178	LEU
1	A	244	GLN

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Mol	Chain	Res	Type
1	A	329	LYS
1	A	347	MET
1	A	375	ASP
1	A	385	ARG
1	A	393	LEU
1	A	427	GLU
1	A	428	ILE
1	B	154	SER
1	B	207	LYS
1	B	244	GLN
1	B	293	LEU
1	B	331	ARG
1	B	360	MET
1	B	375	ASP
1	B	380	ARG
1	B	401	GLU
1	B	404	ARG
1	B	411	MET
1	B	416	GLN
1	B	432	CYS
1	B	436	THR
1	B	474	GLU
1	B	499	CYS
1	B	500	GLN
1	B	511	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	165	HIS
1	A	244	GLN
1	A	382	GLN
1	A	405	GLN
1	A	452	HIS
1	A	463	GLN
1	B	226	HIS
1	B	228	HIS
1	B	244	GLN
1	B	247	HIS
1	B	284	GLN
1	B	288	ASN
1	B	298	ASN

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Mol	Chain	Res	Type
1	B	376	ASN
1	B	386	ASN
1	B	481	GLN
1	B	500	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

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	NVW	A	901	-	34,36,36	0.83	2 (5%)	46,53,53	2.11	8 (17%)
2	NVW	B	901	-	34,36,36	0.77	2 (5%)	46,53,53	1.50	8 (17%)

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	NVW	A	901	-	-	0/17/30/30	0/3/4/4
2	NVW	B	901	-	-	0/17/30/30	0/3/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	901	NVW	C21-N4	2.17	1.38	1.34
2	B	901	NVW	C1-N3	2.27	1.40	1.36
2	A	901	NVW	C1-N3	2.94	1.41	1.36
2	A	901	NVW	C21-N4	3.11	1.39	1.34

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	901	NVW	O1-S1-C11	-7.23	102.88	109.44
2	B	901	NVW	C14-N3-C1	-3.86	118.85	128.96
2	A	901	NVW	C14-N3-C1	-2.47	122.48	128.96
2	B	901	NVW	C10-C9-N1	-2.28	119.82	122.52
2	A	901	NVW	C12-C11-S1	-2.27	106.91	111.61
2	A	901	NVW	C23-C22-N4	2.15	114.23	108.75
2	B	901	NVW	O2-S1-C10	2.19	110.61	108.49
2	B	901	NVW	C10-C1-N3	2.21	122.37	117.99
2	A	901	NVW	C10-C1-N3	2.28	122.52	117.99
2	B	901	NVW	C23-C22-N4	2.47	115.06	108.75
2	B	901	NVW	C12-C13-C9	2.67	117.98	113.48
2	B	901	NVW	O3-C21-N4	2.93	127.66	123.05
2	A	901	NVW	C12-C13-C9	3.76	119.81	113.48
2	A	901	NVW	C13-C9-C10	4.65	124.57	119.13
2	B	901	NVW	C13-C9-C10	5.54	125.61	119.13
2	A	901	NVW	O1-S1-O2	7.42	124.28	117.35

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	NVW	1	0
2	B	901	NVW	1	0

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	345/377 (91%)	0.05	24 (6%)	19	27	28, 43, 78, 110	0
1	B	345/377 (91%)	0.01	20 (5%)	26	35	30, 40, 74, 103	0
All	All	690/754 (91%)	0.03	44 (6%)	23	31	28, 41, 77, 110	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	160	THR	7.5
1	A	157	GLY	5.5
1	A	508	PHE	4.9
1	B	161	GLU	4.6
1	A	511	THR	4.5
1	B	159	ASN	4.5
1	A	256	ASP	4.2
1	A	160	THR	4.1
1	B	157	GLY	4.0
1	A	428	ILE	4.0
1	B	434	LYS	3.3
1	A	161	GLU	3.2
1	A	154	SER	3.2
1	A	165	HIS	3.0
1	A	436	THR	3.0
1	A	432	CYS	2.8
1	B	363	THR	2.8
1	A	243	ALA	2.7
1	A	242	VAL	2.7
1	B	485	PRO	2.6
1	B	436	THR	2.6
1	B	511	THR	2.6
1	B	432	CYS	2.6
1	A	431	MET	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	435	HIS	2.5
1	B	428	ILE	2.5
1	B	359	THR	2.4
1	B	433	ASP	2.4
1	B	242	VAL	2.4
1	A	510	LEU	2.4
1	B	508	PHE	2.4
1	B	371	VAL	2.4
1	A	391	ALA	2.3
1	A	156	PHE	2.3
1	A	158	VAL	2.2
1	A	430	PRO	2.2
1	A	159	ASN	2.2
1	A	239	ALA	2.2
1	B	162	ASN	2.1
1	A	162	ASN	2.1
1	B	360	MET	2.1
1	B	391	ALA	2.1
1	A	501	GLY	2.1
1	A	393	LEU	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	NVW	B	901	33/33	0.96	0.18	0.77	33,43,64,69	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
4	CA	B	903	1/1	1.00	0.16	0.28	38,38,38,38	0
2	NVW	A	901	33/33	0.95	0.14	0.21	42,49,72,73	0
4	CA	A	903	1/1	0.99	0.12	-0.56	40,40,40,40	0
3	ZN	B	902	1/1	1.00	0.13	-1.06	38,38,38,38	0
3	ZN	A	902	1/1	1.00	0.09	-1.62	42,42,42,42	0

6.5 Other polymers (i)

There are no such residues in this entry.