Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 11:32 PM GMT

PDB ID	:	1XOS
Title	:	Catalytic Domain Of Human Phosphodiesterase 4B In Complex With Sildenafil
Authors	:	Card, G.L.; England, B.P.; Suzuki, Y.; Fong, D.; Powell, B.; Lee, B.; Luu, C.; Tabrizizad, M.; Gillette, S.; Ibrahim, P.N.; Artis, D.R.; Bollag, G.; Milburn, M.V.; Kim, S.-H.; Schlessinger, J.; Zhang, K.Y.J.
Deposited on	:	2004-10-06
Resolution	:	2.28 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:

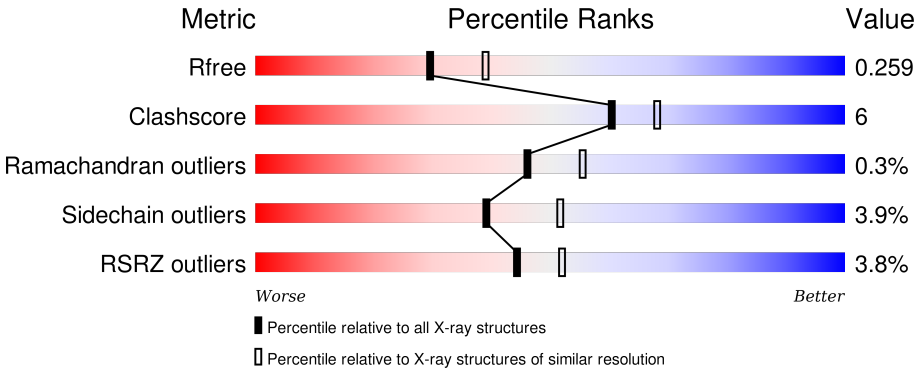
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance

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

The reported resolution of this entry is 2.28 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	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

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	398	3% 67% 13% • 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	ZN	A	1001	-	-	-	X

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 2654 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	322	Total	C	N	O	S	0	0	0
			2597	1641	437	497	22			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	131	MET	-	INITIATING METHIONINE	UNP Q07343
A	132	GLY	-	CLONING ARTIFACT	UNP Q07343
A	133	SER	-	CLONING ARTIFACT	UNP Q07343
A	134	SER	-	CLONING ARTIFACT	UNP Q07343
A	135	HIS	-	EXPRESSION TAG	UNP Q07343
A	136	HIS	-	EXPRESSION TAG	UNP Q07343
A	137	HIS	-	EXPRESSION TAG	UNP Q07343
A	138	HIS	-	EXPRESSION TAG	UNP Q07343
A	139	HIS	-	EXPRESSION TAG	UNP Q07343
A	140	HIS	-	EXPRESSION TAG	UNP Q07343
A	141	SER	-	CLONING ARTIFACT	UNP Q07343
A	142	SER	-	CLONING ARTIFACT	UNP Q07343
A	143	GLY	-	CLONING ARTIFACT	UNP Q07343
A	144	LEU	-	CLONING ARTIFACT	UNP Q07343
A	145	VAL	-	CLONING ARTIFACT	UNP Q07343
A	146	PRO	-	CLONING ARTIFACT	UNP Q07343
A	147	ARG	-	CLONING ARTIFACT	UNP Q07343
A	148	GLY	-	CLONING ARTIFACT	UNP Q07343
A	149	SER	-	CLONING ARTIFACT	UNP Q07343
A	150	HIS	-	CLONING ARTIFACT	UNP Q07343
A	151	MET	-	CLONING ARTIFACT	UNP Q07343
A	194	CME	CYS	MODIFIED RESIDUE	UNP Q07343
A	320	CME	CYS	MODIFIED RESIDUE	UNP Q07343
A	432	CME	CYS	MODIFIED RESIDUE	UNP Q07343

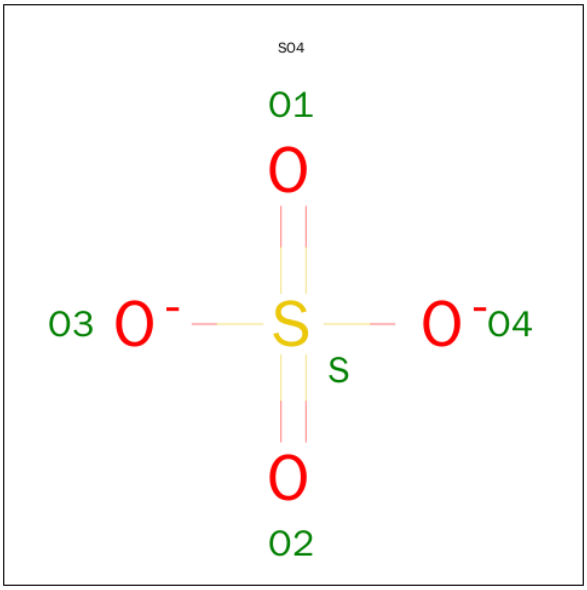
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

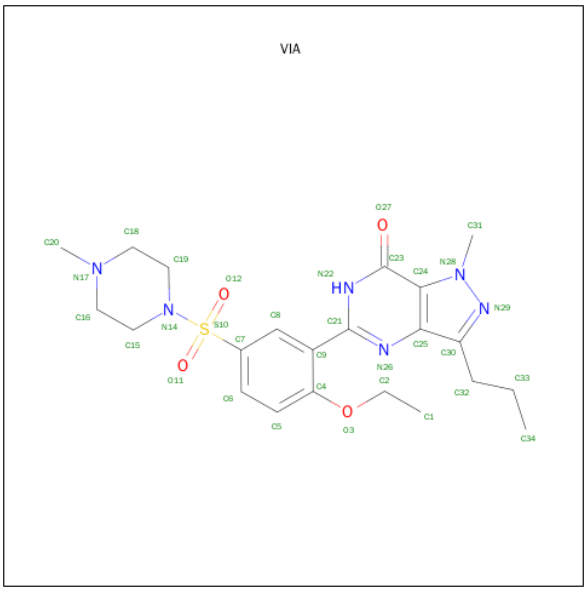
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 5 is 5-{2-ETHOXY-5-[(4-METHYLPIPERAZIN-1-YL)SULFONYL]PHENYL}-1-METHYL-3-PROPYL-1H,6H,7H-PYRAZOLO[4,3-D]PYRIMIDIN-7-ONE (three-letter code: VIA) (formula: $C_{22}H_{30}N_6O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			33	22	6	4	1		

- Molecule 6 is water.

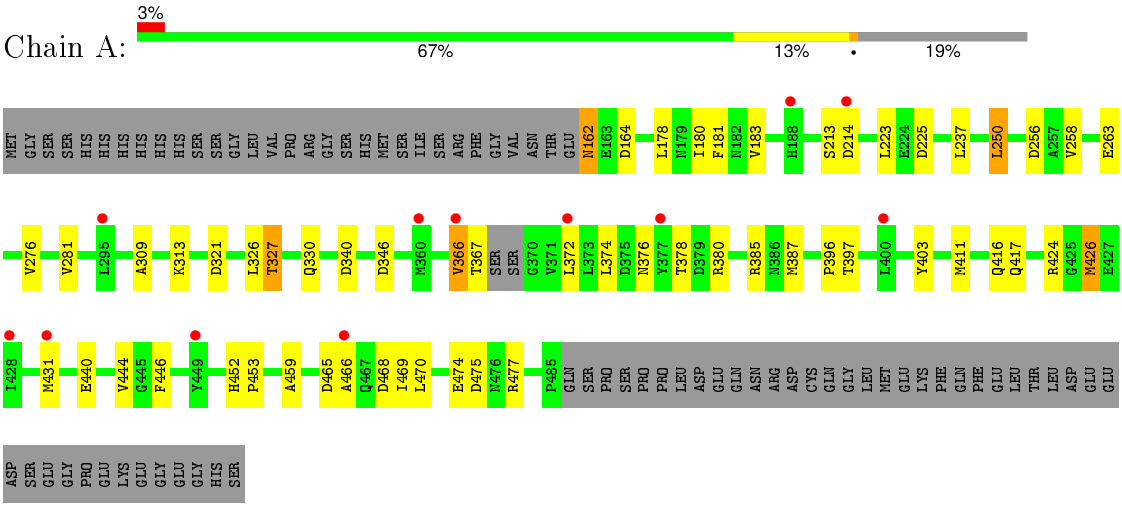
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	17	Total	O	0	0
			17	17		

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:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	94.57Å 107.12Å 89.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	70.71 – 2.28 65.00 – 2.28	Depositor EDS
% Data completeness (in resolution range)	99.7 (70.71-2.28) 99.6 (65.00-2.28)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.17 (at 2.27Å)	Xtriage
Refinement program	REFMAC 5.1.25	Depositor
R, R_{free}	0.210 , 0.257 0.212 , 0.259	Depositor DCC
R_{free} test set	1080 reflections (5.42%)	DCC
Wilson B-factor (Å^2)	52.0	Xtriage
Anisotropy	0.502	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 48.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Outliers	2 of 20997 reflections (0.010%)	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	2654	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 5.80% 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: ZN, MG, VIA, CME, SO4

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	0.68	0/2618	0.87	8/3549 (0.2%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	475	ASP	CB-CG-OD2	8.42	125.88	118.30
1	A	164	ASP	CB-CG-OD2	6.23	123.91	118.30
1	A	256	ASP	CB-CG-OD2	5.70	123.43	118.30
1	A	214	ASP	CB-CG-OD2	5.58	123.32	118.30
1	A	346	ASP	CB-CG-OD1	5.48	123.23	118.30
1	A	225	ASP	CB-CG-OD2	5.24	123.02	118.30
1	A	321	ASP	CB-CG-OD2	5.17	122.96	118.30
1	A	465	ASP	CB-CG-OD2	5.05	122.84	118.30

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	2597	0	2504	28	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	5	0	0	0	0
5	A	33	0	30	5	0
6	A	17	0	0	1	0
All	All	2654	0	2534	29	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 6.

All (29) 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:250:LEU:HD13	1:A:387:MET:CE	2.13	0.78
1:A:250:LEU:HD13	1:A:387:MET:HE3	1.74	0.67
1:A:180:ILE:HG23	1:A:237:LEU:HD11	1.83	0.61
1:A:223:LEU:HD12	1:A:276:VAL:HG21	1.85	0.59
1:A:327:THR:HG22	1:A:330:GLN:H	1.74	0.52
1:A:440:GLU:O	1:A:444:VAL:HG23	2.12	0.50
1:A:396:PRO:HB3	1:A:403:TYR:CZ	2.47	0.49
1:A:424:ARG:HD2	1:A:426:MET:CE	2.42	0.49
1:A:181:PHE:CD2	1:A:237:LEU:HD21	2.47	0.49
1:A:366:VAL:HG23	1:A:367:THR:N	2.29	0.48
1:A:309:ALA:O	1:A:313:LYS:HB2	2.13	0.48
1:A:162:ASN:HD22	1:A:162:ASN:N	2.11	0.48
1:A:397:THR:HB	1:A:469:ILE:HG23	1.95	0.48
1:A:258:VAL:CG1	1:A:374:LEU:HD12	2.43	0.47
1:A:250:LEU:HD13	1:A:387:MET:HE1	1.97	0.47
1:A:180:ILE:HA	1:A:183:VAL:HG13	1.96	0.47
1:A:281:VAL:HG12	1:A:417:GLN:HG3	1.97	0.47
1:A:411:MET:CE	5:A:1:VIA:H321	2.45	0.47
1:A:340:ASP:OD1	1:A:385:ARG:HD3	2.16	0.46
1:A:263:GLU:OE2	1:A:380:ARG:NH1	2.50	0.44
1:A:446:PHE:HB2	5:A:1:VIA:H332	2.00	0.43
1:A:459:ALA:HA	1:A:466:ALA:HB3	2.01	0.43
1:A:470:LEU:O	1:A:474:GLU:HG3	2.19	0.42
1:A:452:HIS:HB3	1:A:453:PRO:HD3	2.01	0.42
1:A:258:VAL:HG11	1:A:374:LEU:HD12	2.02	0.42
1:A:431:MET:SD	5:A:1:VIA:C34	3.08	0.42
5:A:1:VIA:H21	6:A:1008:HOH:O	2.19	0.41
1:A:411:MET:HE3	5:A:1:VIA:H321	2.03	0.41
1:A:376:ASN:HD22	1:A:378:THR:H	1.68	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	315/398 (79%)	301 (96%)	13 (4%)	1 (0%)	46 55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	366	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	283/357 (79%)	272 (96%)	11 (4%)	39 52

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	ASN
1	A	178	LEU
1	A	213	SER
1	A	250	LEU
1	A	326	LEU
1	A	327	THR
1	A	372	LEU

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Mol	Chain	Res	Type
1	A	416	GLN
1	A	426	MET
1	A	468	ASP
1	A	477	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	376	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	CME	A	194	1	8,9,10	1.47	1 (12%)	6,9,11	2.10	3 (50%)
1	CME	A	320	1	8,9,10	1.57	1 (12%)	6,9,11	3.05	2 (33%)
1	CME	A	432	1	8,9,10	1.57	1 (12%)	6,9,11	4.30	3 (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
1	CME	A	194	1	-	0/5/8/10	0/0/0/0
1	CME	A	320	1	-	0/5/8/10	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CME	A	432	1	-	0/5/8/10	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	432	CME	OH-CZ	-3.75	1.21	1.42
1	A	320	CME	OH-CZ	-3.74	1.21	1.42
1	A	194	CME	OH-CZ	-3.58	1.22	1.42

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	320	CME	O-C-CA	-3.61	116.08	125.49
1	A	194	CME	O-C-CA	-2.96	117.78	125.49
1	A	432	CME	O-C-CA	-2.65	118.59	125.49
1	A	432	CME	CZ-CE-SD	-2.55	106.93	113.16
1	A	194	CME	CA-CB-SG	-2.48	105.22	114.23
1	A	194	CME	CB-SG-SD	2.71	109.24	103.95
1	A	320	CME	CB-SG-SD	5.84	115.34	103.95
1	A	432	CME	CB-SG-SD	9.51	122.48	103.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 2 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	$\#\|Z\| > 2$
5	VIA	A	1	-	34,36,36	1.78	7 (20%)	39,53,53	3.05	19 (48%)
4	SO4	A	2	-	4,4,4	0.17	0	6,6,6	0.39	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
5	VIA	A	1	-	-	0/22/32/32	0/4/4/4
4	SO4	A	2	-	-	0/0/0/0	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1	VIA	N29-N28	-3.61	1.33	1.37
5	A	1	VIA	C7-S10	-2.95	1.72	1.76
5	A	1	VIA	C19-N14	-2.29	1.45	1.47
5	A	1	VIA	C30-N29	2.43	1.37	1.33
5	A	1	VIA	C23-C24	3.05	1.47	1.41
5	A	1	VIA	C21-N22	3.68	1.42	1.34
5	A	1	VIA	C23-N22	5.16	1.42	1.33

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	1	VIA	C24-C23-N22	-7.09	113.89	123.59
5	A	1	VIA	C6-C7-C8	-3.41	116.56	120.52
5	A	1	VIA	N26-C21-N22	-2.34	124.69	126.20
5	A	1	VIA	O12-S10-O11	-2.06	115.91	119.47
5	A	1	VIA	C23-C24-C25	-2.04	118.46	120.90
5	A	1	VIA	C9-C21-N22	2.05	120.94	117.43
5	A	1	VIA	C2-O3-C4	2.07	121.88	118.01
5	A	1	VIA	C6-C7-S10	2.11	122.12	119.79
5	A	1	VIA	C19-C18-N17	2.17	113.08	110.79
5	A	1	VIA	C5-C6-C7	2.46	122.20	119.48
5	A	1	VIA	C19-N14-S10	2.47	121.87	117.12
5	A	1	VIA	C20-N17-C18	2.48	114.49	110.63
5	A	1	VIA	C15-N14-S10	2.62	122.15	117.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	1	VIA	C31-N28-N29	2.69	122.53	118.05
5	A	1	VIA	O12-S10-N14	4.20	110.84	106.69
5	A	1	VIA	C23-N22-C21	5.69	121.83	116.23
5	A	1	VIA	C30-N29-N28	5.90	109.51	104.42
5	A	1	VIA	C18-C19-N14	7.48	115.09	109.02
5	A	1	VIA	C16-C15-N14	8.22	115.69	109.02

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1	VIA	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, 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	319/398 (80%)	0.49	12 (3%) 44 52	28, 44, 66, 73	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	372	LEU	3.3
1	A	188	HIS	2.7
1	A	449	TYR	2.6
1	A	466	ALA	2.6
1	A	400	LEU	2.5
1	A	428	ILE	2.4
1	A	366	VAL	2.3
1	A	377	TYR	2.2
1	A	431	MET	2.2
1	A	214	ASP	2.0
1	A	295	LEU	2.0
1	A	360	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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
1	CME	A	432	10/11	0.90	0.20	-	59,62,79,82	0
1	CME	A	194	10/11	0.94	0.14	-	42,45,66,69	0
1	CME	A	320	10/11	0.96	0.14	-	63,66,80,81	0

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	ZN	A	1001	1/1	0.99	0.18	3.17	45,45,45,45	0
5	VIA	A	1	33/33	0.90	0.21	1.43	87,92,95,96	0
3	MG	A	1002	1/1	0.98	0.15	0.65	30,30,30,30	0
4	SO4	A	2	5/5	0.87	0.19	-	109,109,110,110	0

6.5 Other polymers (i)

There are no such residues in this entry.