Full wwPDB X-ray Structure Validation Report (i)

Aug 6, 2023 – 06:51 AM EDT

PDB ID	:	1K57
Title	:	OXA 10 class D beta-lactamase at pH 6.0
Authors	:	Golemi, D.; Maveyraud, L.; Vakulenko, S.; Samama, J.P.; Mobashery, S.
Deposited on	:	2001-10-10
Resolution	:	1.90 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

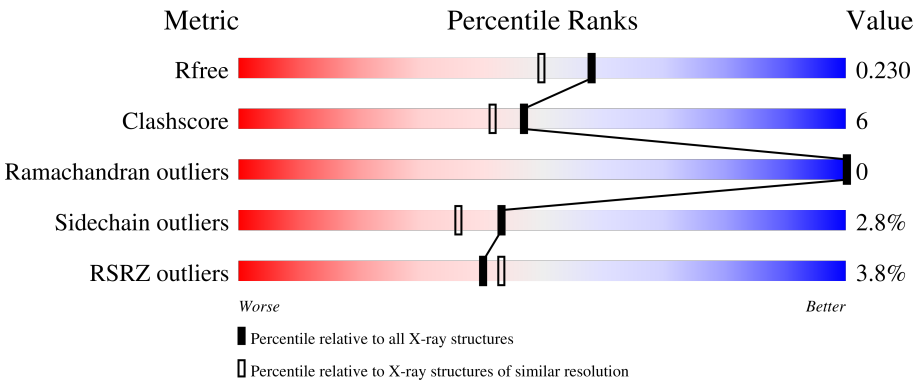
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 1.90 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}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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	246	5% 87% 12%
1	B	246	3% 87% 11%
2	C	246	5% 81% 15%
2	D	246	2% 83% 13%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 8304 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 BETA LACTAMASE OXA-10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	243	1936	1239	324	366	7	0	1	0
1	B	245	1946	1244	326	369	7	0	1	0

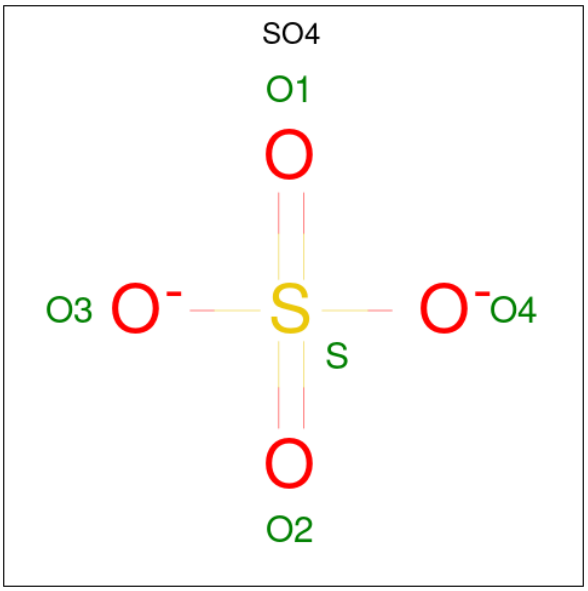
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	70	KCX	LYS	modified residue	UNP P14489
B	70	KCX	LYS	modified residue	UNP P14489

- Molecule 2 is a protein called BETA LACTAMASE OXA-10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	241	1893	1214	314	359	6	0	0	0
2	D	240	1888	1209	313	360	6	0	1	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	118	Total O 118 118	0	0

Continued on next page...

Continued from previous page...

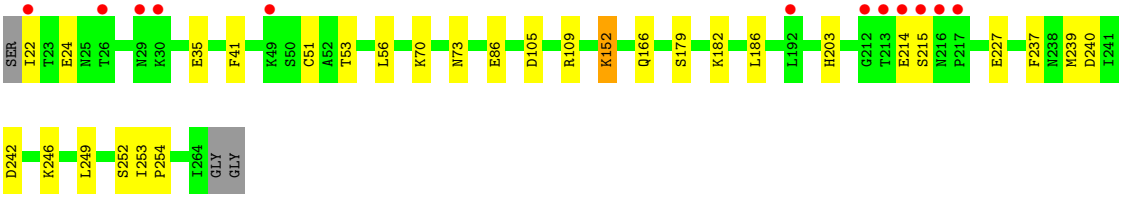
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	B	154	154	154	0	0
			Total	O		
4	C	127	127	127	0	0
			Total	O		
4	D	187	187	187	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

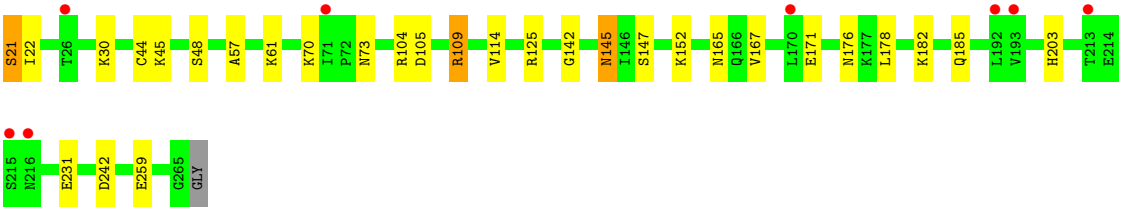
- Molecule 1: BETA LACTAMASE OXA-10

Chain A:



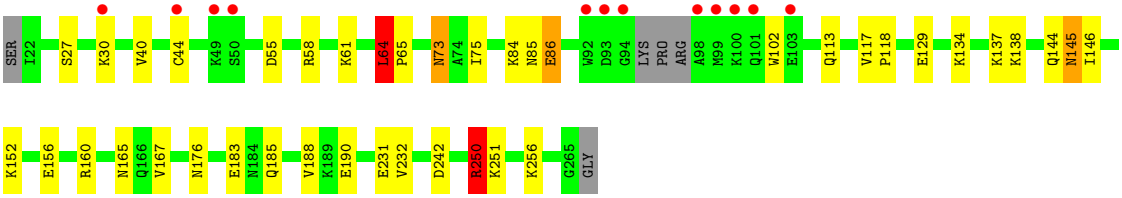
- Molecule 1: BETA LACTAMASE OXA-10

Chain B:



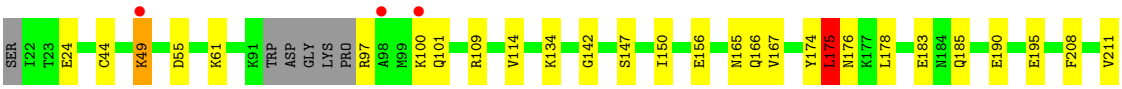
- Molecule 2: BETA LACTAMASE OXA-10

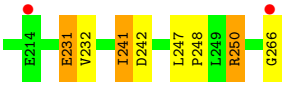
Chain C:



- Molecule 2: BETA LACTAMASE OXA-10

Chain D:




E214
E231
V232
I241
D242
L247
P248
L249
R250
G266

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	66.75Å 82.52Å 101.79Å 90.00° 95.50° 90.00°	Depositor
Resolution (Å)	33.77 – 1.90 33.77 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.6 (33.77-1.90) 97.6 (33.77-1.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	1.88 (at 1.89Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.179 , 0.222 0.190 , 0.230	Depositor DCC
R_{free} test set	1693 reflections (2.00%)	wwPDB-VP
Wilson B-factor (Å^2)	22.2	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 53.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	8304	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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

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	1.00	2/1964 (0.1%)	0.95	4/2653 (0.2%)
1	B	1.11	2/1974 (0.1%)	0.96	5/2666 (0.2%)
2	C	1.08	2/1932 (0.1%)	1.05	8/2611 (0.3%)
2	D	1.13	5/1925 (0.3%)	0.98	4/2599 (0.2%)
All	All	1.08	11/7795 (0.1%)	0.99	21/10529 (0.2%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	250	ARG	CG-CD	-7.40	1.33	1.51
2	D	231	GLU	CD-OE1	6.44	1.32	1.25
1	A	239	MET	SD-CE	-6.33	1.42	1.77
2	C	86	GLU	CD-OE1	6.30	1.32	1.25
2	C	256	LYS	CD-CE	5.99	1.66	1.51
2	D	24	GLU	CD-OE1	5.66	1.31	1.25
2	D	114	VAL	CB-CG2	-5.63	1.41	1.52
1	A	86	GLU	CD-OE2	5.41	1.31	1.25
1	B	259	GLU	CD-OE2	5.29	1.31	1.25
1	B	114	VAL	CB-CG2	-5.21	1.42	1.52
2	D	147	SER	CB-OG	-5.16	1.35	1.42

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	250	ARG	NE-CZ-NH2	-13.20	113.70	120.30
2	C	250	ARG	NE-CZ-NH1	13.10	126.85	120.30
2	C	55	ASP	CB-CG-OD2	8.30	125.77	118.30
2	D	175	LEU	CB-CG-CD1	8.01	124.62	111.00
1	B	242	ASP	CB-CG-OD2	7.85	125.36	118.30
1	A	242	ASP	CB-CG-OD2	7.56	125.11	118.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	105	ASP	CB-CG-OD2	7.17	124.75	118.30
2	C	242	ASP	CB-CG-OD2	6.79	124.41	118.30
2	D	242	ASP	CB-CG-OD2	5.94	123.64	118.30
2	D	55	ASP	CB-CG-OD2	5.79	123.51	118.30
2	D	109	ARG	NE-CZ-NH1	5.77	123.19	120.30
2	C	250	ARG	CD-NE-CZ	5.71	131.59	123.60
2	C	256	LYS	CD-CE-NZ	5.70	124.81	111.70
1	B	109	ARG	NE-CZ-NH2	-5.57	117.52	120.30
2	C	64	LEU	CB-CG-CD1	5.50	120.34	111.00
1	A	109	ARG	NE-CZ-NH2	-5.49	117.56	120.30
1	B	125	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	B	104	ARG	NE-CZ-NH2	-5.41	117.60	120.30
1	A	240	ASP	CB-CG-OD1	5.38	123.14	118.30
1	B	105	ASP	CB-CG-OD2	5.24	123.02	118.30
2	C	64	LEU	CA-CB-CG	5.08	126.98	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1936	0	1928	15	0
1	B	1946	0	1936	27	0
2	C	1893	0	1878	29	0
2	D	1888	0	1878	32	0
3	A	15	0	0	0	0
3	B	20	0	0	0	0
3	C	10	0	0	1	0
3	D	10	0	0	0	0
4	A	118	0	0	3	0
4	B	154	0	0	5	0
4	C	127	0	0	1	0
4	D	187	0	0	6	0
All	All	8304	0	7620	95	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 (95) 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)
2:D:195:GLU:HG3	4:D:1196:HOH:O	1.54	1.06
1:B:182:LYS:HD3	2:D:183:GLU:OE2	1.64	0.98
2:C:84:LYS:HD3	2:C:85:ASN:ND2	1.86	0.89
2:C:134:LYS:HE2	2:C:138:LYS:HE3	1.56	0.85
1:B:22:ILE:N	4:B:1159:HOH:O	2.12	0.82
1:B:21:SER:C	4:B:1159:HOH:O	2.19	0.81
1:B:182:LYS:HD3	2:D:183:GLU:CD	2.01	0.81
2:C:84:LYS:HD3	2:C:85:ASN:HD22	1.47	0.77
2:D:211:VAL:HA	2:D:241[A]:ILE:CD1	2.13	0.77
1:B:182:LYS:CD	2:D:183:GLU:OE2	2.40	0.70
2:C:64:LEU:HD13	2:C:160:ARG:NH2	2.07	0.70
1:B:30:LYS:HD3	1:B:30:LYS:N	2.09	0.66
2:D:266:GLY:HA2	4:D:1170:HOH:O	1.95	0.66
1:B:203:HIS:HD2	4:B:1091:HOH:O	1.80	0.64
2:D:211:VAL:HG22	2:D:241[A]:ILE:HD11	1.78	0.64
1:B:21:SER:HB3	4:B:1121:HOH:O	1.98	0.64
1:B:30:LYS:N	1:B:30:LYS:CD	2.60	0.64
1:A:203:HIS:HD2	4:A:1019:HOH:O	1.82	0.61
1:A:246:LYS:HE2	4:A:1100:HOH:O	1.99	0.61
2:C:134:LYS:CE	2:C:138:LYS:HE3	2.28	0.59
2:D:61:LYS:HE2	4:D:1182:HOH:O	2.03	0.59
2:D:100:LYS:N	2:D:101:GLN:OE1	2.36	0.59
2:D:211:VAL:HA	2:D:241[A]:ILE:HD12	1.83	0.59
2:C:84:LYS:CD	2:C:85:ASN:HD22	2.15	0.58
2:C:250:ARG:NH2	3:C:1004:SO4:O2	2.36	0.58
2:D:101:GLN:H	2:D:101:GLN:CD	2.07	0.58
2:D:208:PHE:HE1	2:D:241[A]:ILE:HG21	1.67	0.57
2:D:49:LYS:HA	2:D:49:LYS:HE2	1.86	0.57
1:B:145:ASN:HD22	1:B:147:SER:H	1.52	0.56
2:C:27:SER:O	2:C:30:LYS:HG2	2.07	0.54
2:C:129:GLU:HG3	2:C:146:ILE:HD11	1.90	0.54
2:D:247:LEU:N	2:D:248:PRO:CD	2.71	0.54
2:C:145:ASN:C	2:C:145:ASN:HD22	2.11	0.53
1:A:186:LEU:HD21	2:C:86:GLU:OE2	2.07	0.53
2:D:241[A]:ILE:C	2:D:241[A]:ILE:HD13	2.30	0.52
2:D:44:CYS:SG	2:D:167:VAL:HG11	2.51	0.51
2:C:61:LYS:HD3	4:C:1134:HOH:O	2.11	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:44:CYS:SG	2:C:167:VAL:HG11	2.51	0.50
2:D:97:ARG:N	4:D:1191:HOH:O	2.44	0.50
2:C:134:LYS:HE2	2:C:138:LYS:CE	2.35	0.49
2:C:176:ASN:HA	2:C:185:GLN:NE2	2.27	0.49
2:C:250:ARG:HG2	2:C:251:LYS:N	2.25	0.49
2:D:150:ILE:O	2:D:150:ILE:HG13	2.11	0.49
1:A:203:HIS:HE1	2:C:190:GLU:OE1	1.95	0.49
2:D:97:ARG:CB	4:D:1139:HOH:O	2.61	0.49
1:B:145:ASN:HD22	1:B:145:ASN:C	2.16	0.49
2:D:208:PHE:CE1	2:D:241[A]:ILE:HG21	2.46	0.49
1:B:145:ASN:ND2	1:B:147:SER:H	2.11	0.49
2:D:178:LEU:H	2:D:185:GLN:HE22	1.60	0.48
1:B:182:LYS:CE	2:D:183:GLU:OE2	2.62	0.48
1:A:182:LYS:HE2	2:C:183:GLU:HG3	1.96	0.48
2:D:100:LYS:HA	2:D:100:LYS:HD2	1.62	0.47
2:C:64:LEU:HD13	2:C:160:ARG:HH21	1.79	0.47
2:C:117:VAL:N	2:C:118:PRO:CD	2.78	0.47
2:C:137:LYS:HA	2:C:144:GLN:NE2	2.30	0.47
2:D:174:TYR:HD2	2:D:175:LEU:HD13	1.80	0.46
1:A:227:GLU:OE1	4:A:1124:HOH:O	2.21	0.46
1:A:22:ILE:HD11	1:A:53:THR:CG2	2.46	0.46
1:B:21:SER:N	4:B:1159:HOH:O	2.49	0.46
1:A:237:PHE:CD1	1:A:237:PHE:C	2.89	0.46
1:B:48:SER:OG	1:B:171:GLU:OE1	2.18	0.45
1:B:176:ASN:ND2	1:B:182:LYS:HE3	2.32	0.45
2:D:174:TYR:CD2	2:D:175:LEU:HD13	2.52	0.45
1:A:152:LYS:HB3	1:A:152:LYS:HE3	1.68	0.45
1:A:249:LEU:HA	1:A:252:SER:OG	2.17	0.44
1:B:142:GLY:H	1:B:165:ASN:HD21	1.64	0.44
1:B:44:CYS:SG	1:B:167:VAL:HG11	2.58	0.44
2:C:40:VAL:HB	2:C:58:ARG:HD3	2.00	0.44
1:B:22:ILE:O	1:B:22:ILE:HG23	2.17	0.43
1:A:41:PHE:HD2	1:A:237:PHE:HB2	1.83	0.43
2:C:134:LYS:CD	2:C:138:LYS:HE3	2.49	0.43
1:A:253:ILE:HB	1:A:254:PRO:HD3	2.00	0.42
1:B:45:LYS:HE3	1:B:231:GLU:HG2	2.01	0.42
1:B:176:ASN:HA	1:B:185:GLN:NE2	2.34	0.42
2:C:113:GLN:HE21	2:C:113:GLN:HB3	1.65	0.42
1:A:22:ILE:HA	1:A:51:CYS:O	2.19	0.42
1:B:203:HIS:HE1	2:D:190:GLU:OE1	2.02	0.42
1:A:35:GLU:HG3	1:A:249:LEU:HD22	2.02	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:ALA:O	1:B:61:LYS:HG3	2.20	0.42
1:A:24:GLU:HB2	1:A:56:LEU:HG	2.00	0.42
1:B:178:LEU:H	1:B:185:GLN:HE22	1.66	0.42
1:B:152:LYS:HB3	1:B:152:LYS:HE2	1.73	0.41
1:B:30:LYS:HD3	1:B:30:LYS:H	1.82	0.41
1:B:109:ARG:HH21	1:B:109:ARG:HD2	1.59	0.41
2:C:75:ILE:HG12	2:C:188:VAL:HG21	2.01	0.41
2:D:142:GLY:H	2:D:165:ASN:ND2	2.19	0.41
2:D:231:GLU:HG3	2:D:232:VAL:N	2.35	0.41
2:C:73:ASN:HD22	2:C:73:ASN:HA	1.68	0.41
2:D:241[A]:ILE:HD13	2:D:241[A]:ILE:O	2.21	0.41
2:C:84:LYS:O	2:C:84:LYS:HG2	2.20	0.40
2:C:231:GLU:HG3	2:C:232:VAL:N	2.36	0.40
2:D:176:ASN:HA	2:D:185:GLN:NE2	2.36	0.40
2:C:102:TRP:CD1	2:C:102:TRP:N	2.88	0.40
2:D:156:GLU:OE1	4:D:1151:HOH:O	2.21	0.40
2:D:211:VAL:HA	2:D:241[A]:ILE:HD13	1.99	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	241/246 (98%)	234 (97%)	7 (3%)	0	100	100
1	B	243/246 (99%)	238 (98%)	5 (2%)	0	100	100
2	C	237/246 (96%)	231 (98%)	6 (2%)	0	100	100
2	D	237/246 (96%)	231 (98%)	6 (2%)	0	100	100
All	All	958/984 (97%)	934 (98%)	24 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	209/209 (100%)	203 (97%)	6 (3%)	42	35
1	B	210/209 (100%)	207 (99%)	3 (1%)	67	65
2	C	204/210 (97%)	196 (96%)	8 (4%)	32	23
2	D	204/210 (97%)	197 (97%)	7 (3%)	37	28
All	All	827/838 (99%)	803 (97%)	24 (3%)	43	35

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	ASN
1	A	152	LYS
1	A	166	GLN
1	A	179	SER
1	A	214	GLU
1	A	215	SER
1	B	21	SER
1	B	73	ASN
1	B	145	ASN
2	C	64	LEU
2	C	65	PRO
2	C	73	ASN
2	C	145	ASN
2	C	152	LYS
2	C	156	GLU
2	C	165	ASN
2	C	250	ARG
2	D	49	LYS
2	D	134	LYS
2	D	166	GLN
2	D	175	LEU
2	D	241[A]	ILE
2	D	241[B]	ILE
2	D	250	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	ASN
1	A	165	ASN
1	A	166	GLN
1	A	203	HIS
1	B	73	ASN
1	B	122	GLN
1	B	145	ASN
1	B	165	ASN
1	B	185	GLN
1	B	203	HIS
2	C	73	ASN
2	C	113	GLN
2	C	121	GLN
2	C	144	GLN
2	C	145	ASN
2	C	185	GLN
2	D	113	GLN
2	D	121	GLN
2	D	165	ASN
2	D	166	GLN
2	D	185	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)

2 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	KCX	A	70	1	9,11,12	1.43	1 (11%)	5,12,14	1.38	1 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
1	KCX	B	70	1	9,11,12	1.21	0	5,12,14	1.64	2 (40%)

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	KCX	A	70	1	-	0/9/10/12	-
1	KCX	B	70	1	-	0/9/10/12	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	70	KCX	CD-CE	2.14	1.60	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	70	KCX	CE-NZ-CX	2.75	126.30	121.89
1	A	70	KCX	OQ1-CX-NZ	-2.46	121.15	124.96
1	B	70	KCX	OQ1-CX-NZ	-2.34	121.32	124.96

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

11 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	$\#\|Z\| > 2$
3	SO4	B	1006	-	4,4,4	0.19	0	6,6,6	0.84	0
3	SO4	B	1003	-	4,4,4	0.54	0	6,6,6	1.14	0
3	SO4	D	1005	-	4,4,4	0.10	0	6,6,6	0.88	0
3	SO4	C	1004	-	4,4,4	0.23	0	6,6,6	0.50	0
3	SO4	D	1010	-	4,4,4	0.34	0	6,6,6	0.90	0
3	SO4	B	1011	-	4,4,4	0.16	0	6,6,6	0.97	0
3	SO4	A	1008	-	4,4,4	0.24	0	6,6,6	0.77	0
3	SO4	B	1007	-	4,4,4	0.11	0	6,6,6	0.36	0
3	SO4	A	1001	-	4,4,4	0.38	0	6,6,6	0.32	0
3	SO4	C	1009	-	4,4,4	0.20	0	6,6,6	0.25	0
3	SO4	A	1002	-	4,4,4	0.29	0	6,6,6	0.80	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1004	SO4	1	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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	242/246 (98%)	0.28	12 (4%)	28	32	14, 23, 39, 52	0
1	B	244/246 (99%)	0.21	8 (3%)	46	49	11, 19, 30, 40	0
2	C	241/246 (97%)	0.23	12 (4%)	28	32	13, 22, 39, 51	0
2	D	240/246 (97%)	0.11	5 (2%)	63	66	9, 18, 33, 45	0
All	All	967/984 (98%)	0.21	37 (3%)	40	43	9, 20, 37, 52	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	216	ASN	5.1
2	C	92	TRP	4.2
2	C	94	GLY	4.0
1	B	215	SER	3.9
1	B	216	ASN	3.8
1	A	213	THR	3.6
2	D	266	GLY	3.6
1	A	217	PRO	3.6
2	C	49	LYS	3.6
2	C	101	GLN	3.5
2	C	30	LYS	3.5
1	A	192	LEU	3.3
1	A	214	GLU	3.2
2	C	100	LYS	3.1
2	C	93	ASP	3.0
1	A	215	SER	2.9
2	C	99	MET	2.9
1	A	26	THR	2.7
2	C	98	ALA	2.7
2	D	100	LYS	2.6
2	D	98	ALA	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	30	LYS	2.5
1	B	213	THR	2.5
2	D	49	LYS	2.5
1	A	212	GLY	2.3
2	C	50	SER	2.3
2	C	103	GLU	2.3
1	B	193	VAL	2.2
1	A	22	ILE	2.2
1	A	49	LYS	2.2
2	C	44	CYS	2.1
1	B	192	LEU	2.1
1	B	71	ILE	2.1
1	A	29	ASN	2.1
1	B	170	LEU	2.1
1	B	26	THR	2.0
2	D	214	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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. 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	B-factors(Å^2)	Q<0.9
1	KCX	A	70	12/13	0.94	0.12	13,15,25,27	0
1	KCX	B	70	12/13	0.94	0.15	10,12,24,26	0

6.3 Carbohydrates (i)

There are no monosaccharides 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. 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	B-factors(Å^2)	Q<0.9
3	SO4	B	1007	5/5	0.90	0.32	61,64,65,65	0
3	SO4	A	1002	5/5	0.91	0.27	35,40,42,46	0
3	SO4	C	1004	5/5	0.93	0.12	38,39,43,45	0
3	SO4	A	1008	5/5	0.94	0.23	40,40,41,41	0
3	SO4	D	1005	5/5	0.95	0.13	30,31,34,41	0
3	SO4	B	1006	5/5	0.97	0.23	42,42,48,49	0
3	SO4	B	1011	5/5	0.97	0.20	29,34,35,36	0
3	SO4	C	1009	5/5	0.98	0.19	30,33,33,35	0
3	SO4	B	1003	5/5	0.98	0.09	25,29,32,33	0
3	SO4	D	1010	5/5	0.98	0.15	24,25,25,26	0
3	SO4	A	1001	5/5	0.99	0.10	26,30,30,33	0

6.5 Other polymers (i)

There are no such residues in this entry.