



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 07:25 PM GMT

PDB ID : 2XWZ
Title : STRUCTURE OF THE RECOMBINANT NATIVE NITRITE REDUCTASE
FROM ALCALIGENES XYLOSOXIDANS COMPLEXED WITH NITRITE
Authors : Antonyuk, S.V.; Leferink, N.G.H.; Han, C.; Heyes, D.J.; Rigby, S.E.J.; Hough,
M.A.; Eady, R.R.; Scrutton, N.S.; Hasnain, S.S.
Deposited on : 2010-11-06
Resolution : 2.34 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

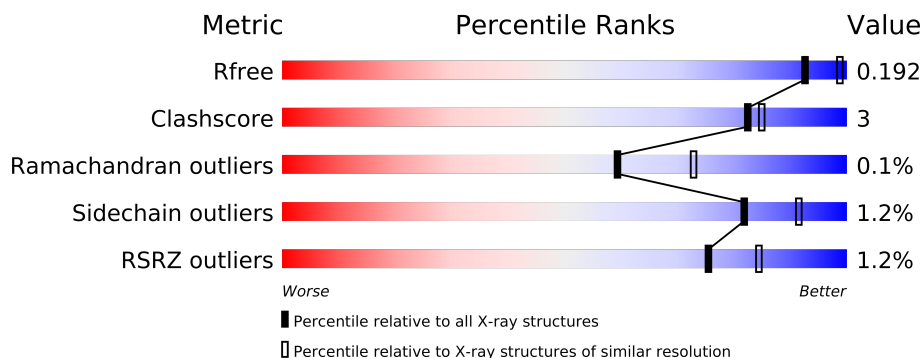
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.34 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}	66092	4049 (2.38-2.30)
Clashscore	79885	1094 (2.36-2.32)
Ramachandran outliers	78287	1080 (2.36-2.32)
Sidechain outliers	78261	1081 (2.36-2.32)
RSRZ outliers	66119	4050 (2.38-2.30)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	336	
1	B	336	
1	C	336	
1	D	336	
1	E	336	
1	F	336	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	NO2	F	1339	-	X
4	SO4	A	1340	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
4	SO4	A	1343	-	X
4	SO4	A	1345	-	X
4	SO4	A	1346	-	X
4	SO4	A	1348	-	X
4	SO4	B	1344	-	X
4	SO4	B	1346	-	X
4	SO4	B	1349	-	X
4	SO4	B	1350	-	X
4	SO4	C	1341	-	X
4	SO4	C	1343	-	X
4	SO4	C	1344	-	X
4	SO4	D	1337	-	X
4	SO4	D	1343	-	X
4	SO4	D	1344	-	X
4	SO4	D	1345	-	X
4	SO4	D	1346[A]	-	X
4	SO4	D	1346[B]	-	X
4	SO4	D	1347	-	X
4	SO4	D	1350	-	X
4	SO4	E	1343	-	X
4	SO4	E	1344	-	X
4	SO4	E	1346	-	X
4	SO4	F	1341	-	X
4	SO4	F	1343	-	X
4	SO4	F	1344	-	X
5	ACT	B	1343	-	X
5	ACT	E	1339	-	X
5	ACT	F	1336	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 17342 atoms, of which 0 are hydrogen and 0 are deuterium.

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 DISSIMILATORY COPPER-CONTAINING NITRITE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	335	Total	C	N	O	S	0	3	0
			2587	1647	450	479	11			
1	B	335	Total	C	N	O	S	0	1	0
			2569	1635	446	477	11			
1	C	335	Total	C	N	O	S	0	2	0
			2576	1640	447	478	11			
1	D	335	Total	C	N	O	S	0	3	0
			2582	1644	448	479	11			
1	E	335	Total	C	N	O	S	0	2	0
			2574	1638	446	479	11			
1	F	334	Total	C	N	O	S	0	5	0
			2586	1648	447	480	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLU	-	EXPRESSION TAG	UNP O68601
B	1	GLU	-	EXPRESSION TAG	UNP O68601
C	1	GLU	-	EXPRESSION TAG	UNP O68601
D	1	GLU	-	EXPRESSION TAG	UNP O68601
E	1	GLU	-	EXPRESSION TAG	UNP O68601
F	1	GLU	-	EXPRESSION TAG	UNP O68601

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

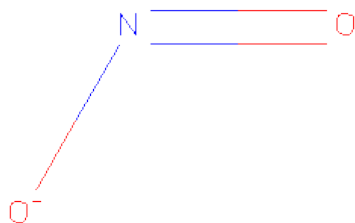
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	2	Total	Cu	0	0
			2	2		
2	E	2	Total	Cu	0	0
			2	2		
2	B	2	Total	Cu	0	0
			2	2		

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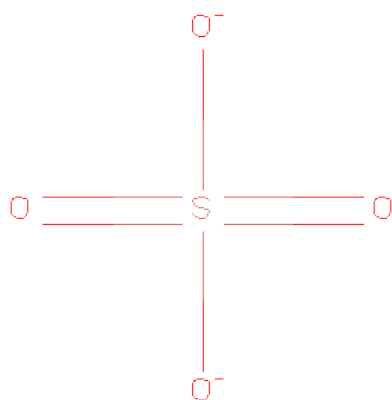
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	2	Total	Cu	0	0
			2	2		
2	A	2	Total	Cu	0	0
			2	2		
2	F	2	Total	Cu	0	0
			2	2		

- Molecule 3 is NITRITE ION (three-letter code: NO₂) (formula: NO₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	N	O	0	0
			3	1	2		
3	A	1	Total	N	O	0	0
			3	1	2		
3	B	1	Total	N	O	0	0
			3	1	2		
3	F	1	Total	N	O	0	0
			3	1	2		

- Molecule 4 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



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

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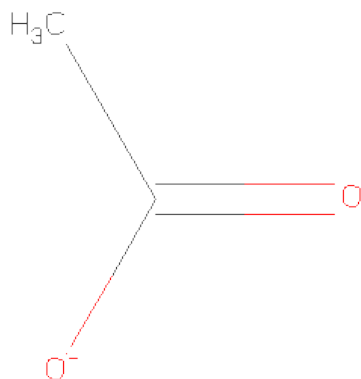
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	1
			10	8	2		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C O 4 2 2	0	0
5	B	1	Total C O 4 2 2	0	0
5	C	1	Total C O 4 2 2	0	0
5	D	1	Total C O 4 2 2	0	0
5	D	1	Total C O 4 2 2	0	0
5	E	1	Total C O 4 2 2	0	0
5	E	1	Total C O 4 2 2	0	0
5	E	1	Total C O 4 2 2	0	0
5	F	1	Total C O 4 2 2	0	0

- Molecule 6 is NITRIC OXIDE (three-letter code: NO) (formula: NO).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	C	1	Total N O 2 1 1	0	0
6	E	1	Total N O 2 1 1	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	275	Total 275	O 275	0	0
7	B	234	Total 234	O 234	0	0
7	C	258	Total 258	O 258	0	0
7	D	261	Total 261	O 261	0	0
7	E	277	Total 277	O 277	0	0
7	F	266	Total 266	O 266	0	0

3 Residue-property plots

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

- Molecule 1: DISSIMILATORY COPPER-CONTAINING NITRITE REDUCTASE

Chain A: 



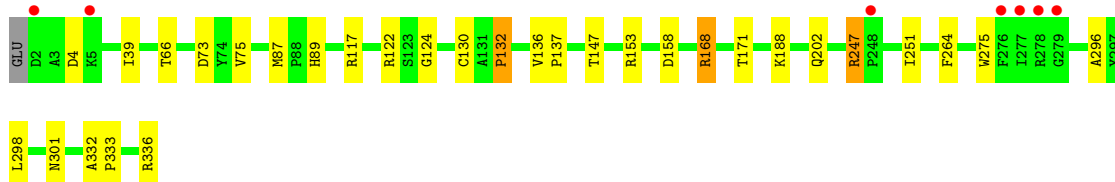
- Molecule 1: DISSIMILATORY COPPER-CONTAINING NITRITE REDUCTASE

Chain B: 



- Molecule 1: DISSIMILATORY COPPER-CONTAINING NITRITE REDUCTASE

Chain C: 



- Molecule 1: DISSIMILATORY COPPER-CONTAINING NITRITE REDUCTASE

Chain D: 



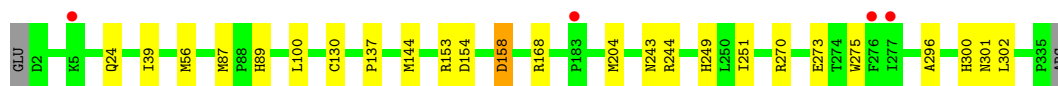
- Molecule 1: DISSIMILATORY COPPER-CONTAINING NITRITE REDUCTASE

Chain E: 



- Molecule 1: DISSIMILATORY COPPER-CONTAINING NITRITE REDUCTASE

Chain F: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	173.93Å 176.82Å 181.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.97 – 2.34 43.97 – 2.34	Depositor EDS
% Data completeness (in resolution range)	99.7 (43.97-2.34) 99.4 (43.97-2.34)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.174 , 0.199 0.168 , 0.192	Depositor DCC
R_{free} test set	11648 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	46.7	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 31.7	EDS
Estimated twinning fraction	0.004 for -h,l,k 0.004 for -l,-k,-h 0.007 for k,h,-l 0.000 for k,l,h 0.000 for l,h,k	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 232959 reflections (0.000%)	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	17342	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.52% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, ACT, NO2, CU, NO

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	0.75	0/2667	0.78	3/3629 (0.1%)
1	B	0.73	0/2643	0.77	2/3599 (0.1%)
1	C	0.78	0/2653	0.78	1/3612 (0.0%)
1	D	0.77	0/2662	0.78	1/3624 (0.0%)
1	E	0.78	0/2651	0.76	1/3610 (0.0%)
1	F	0.76	1/2672 (0.0%)	0.77	2/3638 (0.1%)
All	All	0.76	1/15948 (0.0%)	0.78	10/21712 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	D	0	1
1	E	0	1
1	F	0	1
All	All	0	5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	130	CYS	CB-SG	-5.56	1.72	1.81

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	153	ARG	NE-CZ-NH2	-6.26	117.17	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	153	ARG	CG-CD-NE	-5.83	99.56	111.80
1	A	144	MET	CG-SD-CE	-5.78	90.96	100.20
1	B	158	ASP	CB-CG-OD1	5.73	123.46	118.30
1	B	247	ARG	NE-CZ-NH1	5.27	122.94	120.30
1	F	244	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	E	135	MET	CG-SD-CE	5.12	108.39	100.20
1	D	153	ARG	NE-CZ-NH2	-5.11	117.75	120.30
1	F	153	ARG	NE-CZ-NH2	-5.08	117.76	120.30
1	C	122	ARG	NE-CZ-NH1	5.04	122.82	120.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	300	HIS	Peptide
1	B	300	HIS	Peptide
1	D	300	HIS	Peptide
1	E	300	HIS	Peptide
1	F	300	HIS	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2587	0	2555	15	0
1	B	2569	0	2525	16	0
1	C	2576	0	2535	21	0
1	D	2582	0	2539	23	0
1	E	2574	0	2529	10	0
1	F	2586	0	2554	17	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	2	0	0	0	0
3	A	6	0	0	0	0
3	B	3	0	0	0	0
3	F	3	0	0	0	0
4	A	35	0	0	0	0
4	B	40	0	0	2	0
4	C	30	0	0	0	0
4	D	50	0	0	1	0
4	E	35	0	0	0	0
4	F	35	0	0	2	0
5	A	4	0	3	0	0
5	B	12	0	9	4	0
5	C	4	0	3	4	0
5	D	8	0	6	2	0
5	E	12	0	9	6	0
5	F	4	0	3	0	0
6	C	2	0	0	0	0
6	E	2	0	0	0	0
7	A	275	0	0	0	0
7	B	234	0	0	6	0
7	C	258	0	0	3	0
7	D	261	0	0	3	0
7	E	277	0	0	2	0
7	F	266	0	0	1	0
All	All	17342	0	15270	95	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 3.

All (95) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:87:MET:HE2	7:B:2096:HOH:O	1.59	1.01
1:A:307:GLU:OE2	5:D:1338:ACT:H1	1.68	0.93
1:D:87:MET:HE2	7:D:2116:HOH:O	1.75	0.85
1:B:87:MET:CE	7:B:2096:HOH:O	2.17	0.84
1:D:290:LYS:NZ	5:E:1339:ACT:H1	1.96	0.81
1:D:202[B]:GLN:OE1	1:D:205[B]:ARG:NH1	2.15	0.80
1:B:251:ILE:HD12	1:B:296:ALA:HB3	1.69	0.74
4:B:1345:SO4:O1	7:B:2229:HOH:O	2.07	0.72
1:C:202[B]:GLN:OE1	7:C:2168:HOH:O	2.08	0.71
4:B:1347:SO4:O4	7:B:2230:HOH:O	2.07	0.70
1:B:307:GLU:OE2	5:C:1337:ACT:H3	1.93	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:39:ILE:HD11	1:B:79:LEU:HD11	1.75	0.68
1:D:290:LYS:HZ1	5:E:1339:ACT:H1	1.56	0.68
1:E:188:LYS:HD2	7:E:2172:HOH:O	1.95	0.66
1:E:251:ILE:HD12	1:E:296:ALA:HB3	1.78	0.65
1:B:307:GLU:OE2	5:C:1337:ACT:CH3	2.44	0.65
4:D:1348:SO4:O3	7:D:2260:HOH:O	2.12	0.65
1:C:188:LYS:HD2	7:C:2166:HOH:O	1.98	0.63
1:B:137:PRO:HB2	1:B:204:MET:HE1	1.80	0.63
1:A:20:HIS:CE1	1:A:68:VAL:H	2.18	0.62
5:B:1343:ACT:H1	1:F:270:ARG:NH2	2.15	0.61
1:F:154[B]:ASP:HB2	4:F:1342:SO4:O2	2.01	0.61
1:D:137:PRO:HG2	5:D:1338:ACT:H2	1.81	0.61
1:D:137:PRO:HB2	1:D:204:MET:HE1	1.84	0.60
1:D:290:LYS:HZ3	5:E:1339:ACT:H1	1.69	0.57
1:A:100:LEU:HA	1:E:251:ILE:HG22	1.87	0.56
1:D:251:ILE:HD12	1:D:296:ALA:HB3	1.89	0.54
1:C:168:ARG:NH2	7:C:2131:HOH:O	2.40	0.54
1:A:20:HIS:HE1	1:A:68:VAL:H	1.56	0.53
5:B:1343:ACT:H2	7:B:2058:HOH:O	2.09	0.53
1:C:87:MET:HA	1:C:87:MET:HE3	1.91	0.53
1:D:39:ILE:HD13	1:D:89:HIS:HB2	1.93	0.51
1:C:251:ILE:HG22	1:F:100:LEU:HA	1.93	0.51
1:F:87:MET:HE2	7:F:2126:HOH:O	2.09	0.51
1:F:137:PRO:HB2	1:F:204:MET:HE1	1.92	0.51
1:C:336:ARG:N	1:C:336:ARG:HD2	2.28	0.49
1:B:136:VAL:HG12	1:F:302:LEU:HD13	1.94	0.48
1:A:251:ILE:HD12	1:A:296:ALA:HB3	1.95	0.48
1:B:100:LEU:HA	1:F:251:ILE:HG22	1.96	0.48
1:A:39:ILE:HD13	1:A:89:HIS:HB2	1.96	0.47
1:D:20:HIS:CE1	1:D:68:VAL:H	2.33	0.47
1:E:134:GLY:HA2	5:E:1338:ACT:H3	1.95	0.47
1:D:42:LYS:NZ	1:D:60:GLY:O	2.40	0.47
1:B:171:THR:O	1:B:171:THR:HG23	2.15	0.47
1:D:251:ILE:HG22	1:E:100:LEU:HA	1.96	0.47
5:B:1343:ACT:CH3	7:B:2058:HOH:O	2.62	0.46
1:C:73:ASP:OD1	1:C:153:ARG:NH2	2.42	0.46
1:D:247:ARG:HA	1:D:275:TRP:O	2.15	0.46
1:D:20:HIS:HE1	1:D:68:VAL:H	1.64	0.46
1:C:66:THR:HA	1:C:147:THR:O	2.16	0.46
1:A:301:ASN:HA	1:D:243:ASN:O	2.15	0.46
1:C:137:PRO:HB2	5:C:1337:ACT:H1	1.98	0.46
1:B:42:LYS:NZ	1:B:60:GLY:O	2.46	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:251:ILE:HD12	1:C:296:ALA:HB3	1.97	0.46
1:F:249:HIS:ND1	1:F:273:GLU:O	2.49	0.45
1:A:217:LYS:HG2	1:A:220:ALA:HB2	1.98	0.45
1:D:137:PRO:HB2	1:D:204:MET:CE	2.46	0.45
1:C:301:ASN:HA	1:F:243:ASN:O	2.16	0.45
1:A:294:VAL:HG22	1:A:315:LYS:HG3	1.98	0.44
1:C:332:ALA:HB1	1:C:333:PRO:HD2	1.98	0.44
1:C:336:ARG:NH2	5:E:1338:ACT:O	2.50	0.44
1:E:73:ASP:OD1	1:E:153:ARG:NH2	2.35	0.44
1:B:299:ASN:O	1:B:305:ALA:HB2	2.17	0.44
1:A:136:VAL:HB	1:A:137:PRO:HD3	1.99	0.43
1:F:137:PRO:HB2	1:F:204:MET:CE	2.48	0.43
1:A:276:PHE:HB2	1:D:278:ARG:HG2	2.00	0.43
1:E:202[B]:GLN:NE2	7:E:2178:HOH:O	2.41	0.43
1:B:243:ASN:O	1:F:301:ASN:HA	2.19	0.43
1:D:66:THR:HA	1:D:147:THR:O	2.19	0.43
1:C:75:VAL:O	1:C:117:ARG:HA	2.19	0.43
1:A:249:HIS:ND1	1:A:273:GLU:O	2.52	0.43
5:B:1343:ACT:H1	1:F:270:ARG:HH22	1.84	0.43
1:A:130:CYS:SG	1:A:132:PRO:HD3	2.59	0.42
1:F:154[B]:ASP:CB	4:F:1342:SO4:O2	2.68	0.42
1:E:257:TRP:CZ3	1:E:270:ARG:HB3	2.54	0.42
1:D:307:GLU:OE2	5:E:1337:ACT:O	2.38	0.42
1:D:298:LEU:C	1:D:298:LEU:HD12	2.40	0.42
1:A:137:PRO:HB2	1:A:204:MET:CE	2.50	0.42
1:F:56:MET:HB2	1:F:144:MET:HG3	2.01	0.42
1:D:336:ARG:NH1	7:D:2249:HOH:O	2.53	0.42
1:B:251:ILE:CD1	1:B:296:ALA:HB3	2.46	0.41
1:D:32:VAL:HG22	1:D:74:TYR:HB2	2.02	0.41
1:C:39:ILE:HD13	1:C:89:HIS:HB2	2.02	0.41
1:C:298:LEU:C	1:C:298:LEU:HD12	2.41	0.41
1:F:251:ILE:HD12	1:F:296:ALA:HB3	2.02	0.41
1:F:39:ILE:HD13	1:F:89:HIS:HB2	2.02	0.41
1:C:130:CYS:SG	1:C:132:PRO:HD3	2.61	0.41
1:C:247:ARG:HA	1:C:275:TRP:O	2.21	0.41
1:F:158:ASP:C	1:F:158:ASP:OD1	2.59	0.40
1:C:171:THR:O	1:C:171:THR:HG23	2.21	0.40
1:C:124:GLY:HA2	1:C:264:PHE:CD2	2.56	0.40
1:A:243:ASN:O	1:E:301:ASN:HA	2.21	0.40
1:B:302:LEU:HD13	1:C:136:VAL:HG12	2.03	0.40
1:B:307:GLU:OE2	5:C:1337:ACT:H2	2.19	0.40
1:D:301:ASN:HA	1:E:243:ASN:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	336/336 (100%)	328 (98%)	8 (2%)	0	100	100
1	B	334/336 (99%)	326 (98%)	7 (2%)	1 (0%)	50	60
1	C	335/336 (100%)	325 (97%)	10 (3%)	0	100	100
1	D	336/336 (100%)	324 (96%)	12 (4%)	0	100	100
1	E	335/336 (100%)	326 (97%)	9 (3%)	0	100	100
1	F	337/336 (100%)	328 (97%)	8 (2%)	1 (0%)	50	60
All	All	2013/2016 (100%)	1957 (97%)	54 (3%)	2 (0%)	59	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	24	GLN
1	B	24	GLN

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	273/271 (101%)	271 (99%)	2 (1%)	91	96
1	B	270/271 (100%)	266 (98%)	4 (2%)	76	88
1	C	271/271 (100%)	266 (98%)	5 (2%)	71	84
1	D	271/271 (100%)	268 (99%)	3 (1%)	84	92
1	E	271/271 (100%)	268 (99%)	3 (1%)	84	92

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	274/271 (101%)	271 (99%)	3 (1%)	84	92
All	All	1630/1626 (100%)	1610 (99%)	20 (1%)	82	91

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

Mol	Chain	Res	Type
1	A	158	ASP
1	A	336	ARG
1	B	9	THR
1	B	168	ARG
1	B	247	ARG
1	B	336	ARG
1	C	4	ASP
1	C	132	PRO
1	C	158	ASP
1	C	168	ARG
1	C	247	ARG
1	D	133	GLU
1	D	158	ASP
1	D	336	ARG
1	E	135	MET
1	E	158	ASP
1	E	336	ARG
1	F	158	ASP
1	F	168	ARG
1	F	275	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	20	HIS
1	D	20	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 74 ligands modelled in this entry, 12 are monoatomic - leaving 62 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	$\# Z > 2$
3	NO2	A	1339	2	2,2,2	1.71	0	1,1,1	0.32	0
4	SO4	A	1340	-	4,4,4	0.15	0	6,6,6	0.43	0
5	ACT	A	1341	-	1,3,3	2.53	1 (100%)	0,3,3	0.00	-
3	NO2	A	1342	2	2,2,2	1.64	0	1,1,1	0.35	0
4	SO4	A	1343	-	4,4,4	0.44	0	6,6,6	0.43	0
4	SO4	A	1344	-	4,4,4	0.38	0	6,6,6	0.60	0
4	SO4	A	1345	-	4,4,4	0.09	0	6,6,6	0.25	0
4	SO4	A	1346	-	4,4,4	0.27	0	6,6,6	0.53	0
4	SO4	A	1347	-	4,4,4	0.13	0	6,6,6	0.48	0
4	SO4	A	1348	-	4,4,4	0.14	0	6,6,6	0.16	0
5	ACT	B	1337	-	1,3,3	1.51	0	0,3,3	0.00	-
4	SO4	B	1340	-	4,4,4	0.21	0	6,6,6	0.31	0
3	NO2	B	1341	2	2,2,2	1.64	0	1,1,1	0.29	0
5	ACT	B	1342	-	1,3,3	2.10	1 (100%)	0,3,3	0.00	-
5	ACT	B	1343	-	1,3,3	0.58	0	0,3,3	0.00	-
4	SO4	B	1344	-	4,4,4	0.44	0	6,6,6	0.27	0
4	SO4	B	1345	-	4,4,4	0.30	0	6,6,6	0.46	0
4	SO4	B	1346	-	4,4,4	0.41	0	6,6,6	0.70	0
4	SO4	B	1347	-	4,4,4	0.11	0	6,6,6	0.17	0
4	SO4	B	1348	-	4,4,4	0.10	0	6,6,6	0.19	0
4	SO4	B	1349	-	4,4,4	0.15	0	6,6,6	0.28	0
4	SO4	B	1350	-	4,4,4	0.27	0	6,6,6	0.58	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	ACT	C	1337	-	1,3,3	0.66	0	0,3,3	0.00	-
4	SO4	C	1341	-	4,4,4	0.21	0	6,6,6	0.25	0
4	SO4	C	1342	-	4,4,4	0.24	0	6,6,6	0.46	0
4	SO4	C	1343	-	4,4,4	0.32	0	6,6,6	0.39	0
4	SO4	C	1344	-	4,4,4	0.37	0	6,6,6	0.47	0
4	SO4	C	1345	-	4,4,4	0.25	0	6,6,6	0.83	0
4	SO4	C	1346	-	4,4,4	0.42	0	6,6,6	0.34	0
6	NO	C	1347	2	0,1,1	0.00	-	0,0,0	0.00	-
4	SO4	D	1337	-	4,4,4	0.26	0	6,6,6	0.56	0
5	ACT	D	1338	-	1,3,3	1.24	0	0,3,3	0.00	-
5	ACT	D	1339	-	1,3,3	1.42	0	0,3,3	0.00	-
4	SO4	D	1343	-	4,4,4	0.29	0	6,6,6	0.60	0
4	SO4	D	1344	-	4,4,4	0.21	0	6,6,6	0.73	0
4	SO4	D	1345	-	4,4,4	0.22	0	6,6,6	0.59	0
4	SO4	D	1346[A]	-	4,4,4	0.20	0	6,6,6	0.26	0
4	SO4	D	1346[B]	-	4,4,4	0.21	0	6,6,6	0.54	0
4	SO4	D	1347	-	4,4,4	0.22	0	6,6,6	0.24	0
4	SO4	D	1348	-	4,4,4	0.26	0	6,6,6	0.46	0
4	SO4	D	1349	-	4,4,4	0.16	0	6,6,6	0.37	0
4	SO4	D	1350	-	4,4,4	0.23	0	6,6,6	0.46	0
5	ACT	E	1337	-	1,3,3	2.38	1 (100%)	0,3,3	0.00	-
5	ACT	E	1338	-	1,3,3	0.51	0	0,3,3	0.00	-
5	ACT	E	1339	-	1,3,3	0.93	0	0,3,3	0.00	-
4	SO4	E	1342	-	4,4,4	0.48	0	6,6,6	0.42	0
4	SO4	E	1343	-	4,4,4	0.09	0	6,6,6	0.41	0
4	SO4	E	1344	-	4,4,4	0.29	0	6,6,6	0.16	0
4	SO4	E	1345	-	4,4,4	0.20	0	6,6,6	0.34	0
4	SO4	E	1346	-	4,4,4	0.24	0	6,6,6	0.43	0
4	SO4	E	1347	-	4,4,4	0.22	0	6,6,6	0.46	0
4	SO4	E	1348	-	4,4,4	0.35	0	6,6,6	0.52	0
6	NO	E	1349	2	0,1,1	0.00	-	0,0,0	0.00	-
5	ACT	F	1336	-	1,3,3	3.71	1 (100%)	0,3,3	0.00	-
3	NO2	F	1339	2	2,2,2	1.75	0	1,1,1	0.35	0
4	SO4	F	1340	-	4,4,4	0.39	0	6,6,6	0.28	0
4	SO4	F	1341	-	4,4,4	0.14	0	6,6,6	0.53	0
4	SO4	F	1342	-	4,4,4	0.23	0	6,6,6	0.34	0
4	SO4	F	1343	-	4,4,4	0.21	0	6,6,6	0.35	0
4	SO4	F	1344	-	4,4,4	0.44	0	6,6,6	0.42	0
4	SO4	F	1345	-	4,4,4	0.49	0	6,6,6	0.37	0
4	SO4	F	1346	-	4,4,4	0.23	0	6,6,6	0.38	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
3	NO2	A	1339	2	-	0/0/0/0	0/0/0/0
4	SO4	A	1340	-	-	0/0/0/0	0/0/0/0
5	ACT	A	1341	-	-	0/0/0/0	0/0/0/0
3	NO2	A	1342	2	-	0/0/0/0	0/0/0/0
4	SO4	A	1343	-	-	0/0/0/0	0/0/0/0
4	SO4	A	1344	-	-	0/0/0/0	0/0/0/0
4	SO4	A	1345	-	-	0/0/0/0	0/0/0/0
4	SO4	A	1346	-	-	0/0/0/0	0/0/0/0
4	SO4	A	1347	-	-	0/0/0/0	0/0/0/0
4	SO4	A	1348	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1337	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1340	-	-	0/0/0/0	0/0/0/0
3	NO2	B	1341	2	-	0/0/0/0	0/0/0/0
5	ACT	B	1342	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1343	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1344	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1345	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1346	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1347	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1348	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1349	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1350	-	-	0/0/0/0	0/0/0/0
5	ACT	C	1337	-	-	0/0/0/0	0/0/0/0
4	SO4	C	1341	-	-	0/0/0/0	0/0/0/0
4	SO4	C	1342	-	-	0/0/0/0	0/0/0/0
4	SO4	C	1343	-	-	0/0/0/0	0/0/0/0
4	SO4	C	1344	-	-	0/0/0/0	0/0/0/0
4	SO4	C	1345	-	-	0/0/0/0	0/0/0/0
4	SO4	C	1346	-	-	0/0/0/0	0/0/0/0
6	NO	C	1347	2	-	0/0/0/0	0/0/0/0
4	SO4	D	1337	-	-	0/0/0/0	0/0/0/0
5	ACT	D	1338	-	-	0/0/0/0	0/0/0/0
5	ACT	D	1339	-	-	0/0/0/0	0/0/0/0
4	SO4	D	1343	-	-	0/0/0/0	0/0/0/0
4	SO4	D	1344	-	-	0/0/0/0	0/0/0/0
4	SO4	D	1345	-	-	0/0/0/0	0/0/0/0
4	SO4	D	1346[A]	-	-	0/0/0/0	0/0/0/0
4	SO4	D	1346[B]	-	-	0/0/0/0	0/0/0/0
4	SO4	D	1347	-	-	0/0/0/0	0/0/0/0
4	SO4	D	1348	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SO4	D	1349	-	-	0/0/0/0	0/0/0/0
4	SO4	D	1350	-	-	0/0/0/0	0/0/0/0
5	ACT	E	1337	-	-	0/0/0/0	0/0/0/0
5	ACT	E	1338	-	-	0/0/0/0	0/0/0/0
5	ACT	E	1339	-	-	0/0/0/0	0/0/0/0
4	SO4	E	1342	-	-	0/0/0/0	0/0/0/0
4	SO4	E	1343	-	-	0/0/0/0	0/0/0/0
4	SO4	E	1344	-	-	0/0/0/0	0/0/0/0
4	SO4	E	1345	-	-	0/0/0/0	0/0/0/0
4	SO4	E	1346	-	-	0/0/0/0	0/0/0/0
4	SO4	E	1347	-	-	0/0/0/0	0/0/0/0
4	SO4	E	1348	-	-	0/0/0/0	0/0/0/0
6	NO	E	1349	2	-	0/0/0/0	0/0/0/0
5	ACT	F	1336	-	-	0/0/0/0	0/0/0/0
3	NO2	F	1339	2	-	0/0/0/0	0/0/0/0
4	SO4	F	1340	-	-	0/0/0/0	0/0/0/0
4	SO4	F	1341	-	-	0/0/0/0	0/0/0/0
4	SO4	F	1342	-	-	0/0/0/0	0/0/0/0
4	SO4	F	1343	-	-	0/0/0/0	0/0/0/0
4	SO4	F	1344	-	-	0/0/0/0	0/0/0/0
4	SO4	F	1345	-	-	0/0/0/0	0/0/0/0
4	SO4	F	1346	-	-	0/0/0/0	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	1336	ACT	CH3-C	3.71	1.54	1.48
5	A	1341	ACT	CH3-C	2.53	1.52	1.48
5	E	1337	ACT	CH3-C	2.38	1.52	1.48
5	B	1342	ACT	CH3-C	2.10	1.51	1.48

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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, 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(Å²)	Q<0.9	
1	A	335/336 (99%)	-0.24	5 (1%)	70	80	28, 37, 54, 74	3 (0%)
1	B	335/336 (99%)	-0.32	4 (1%)	75	84	28, 39, 58, 85	7 (2%)
1	C	335/336 (99%)	-0.28	7 (2%)	60	71	27, 38, 55, 81	4 (1%)
1	D	335/336 (99%)	-0.36	2 (0%)	86	93	28, 39, 57, 77	5 (1%)
1	E	335/336 (99%)	-0.30	4 (1%)	75	84	28, 38, 54, 77	4 (1%)
1	F	334/336 (99%)	-0.36	4 (1%)	75	84	27, 38, 57, 77	4 (1%)
All	All	2009/2016 (99%)	-0.31	26 (1%)	75	83	27, 38, 56, 85	27 (1%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	276	PHE	3.5
1	E	277	ILE	3.5
1	A	277	ILE	3.1
1	A	276	PHE	3.0
1	C	276	PHE	2.9
1	A	6	LEU	2.8
1	F	183	PRO	2.8
1	D	276	PHE	2.7
1	F	276	PHE	2.7
1	C	277	ILE	2.7
1	B	2	ASP	2.7
1	B	276	PHE	2.6
1	A	5	LYS	2.5
1	E	4	ASP	2.5
1	D	277	ILE	2.4
1	A	278	ARG	2.4
1	E	2	ASP	2.3
1	F	277	ILE	2.3
1	C	2	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	246	THR	2.1
1	C	279	GLY	2.1
1	B	3	ALA	2.1
1	C	5	LYS	2.1
1	C	248	PRO	2.1
1	F	5	LYS	2.1
1	C	278	ARG	2.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	SO4	E	1346	5/5	0.25	17.68	75,75,78,78	5
4	SO4	F	1344	5/5	0.19	14.81	64,66,66,67	5
4	SO4	B	1346	5/5	0.22	12.20	42,45,47,47	5
5	ACT	B	1343	4/4	0.28	10.72	64,64,65,65	0
4	SO4	B	1344	5/5	0.15	10.42	42,44,46,47	5
4	SO4	A	1346	5/5	0.23	7.93	61,62,63,65	5
4	SO4	E	1344	5/5	0.12	7.37	42,44,45,46	5
4	SO4	A	1348	5/5	0.34	7.04	70,70,71,71	5
4	SO4	D	1344	5/5	0.17	6.74	47,49,50,51	5
4	SO4	D	1346[A]	5/5	0.19	6.71	49,50,51,52	5
4	SO4	F	1343	5/5	0.19	6.39	41,45,46,46	5
4	SO4	B	1350	5/5	0.18	6.26	49,50,51,51	5
4	SO4	F	1341	5/5	0.21	5.88	48,50,53,53	5
4	SO4	D	1347	5/5	0.20	5.73	63,64,65,65	5

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	SO4	D	1345	5/5	0.15	5.07	42,44,45,45	5
4	SO4	D	1346[B]	5/5	0.19	4.44	51,52,57,58	5
4	SO4	C	1341	5/5	0.12	4.33	55,56,57,57	5
4	SO4	C	1344	5/5	0.18	4.26	46,48,51,53	5
4	SO4	A	1343	5/5	0.12	4.25	49,49,50,51	5
4	SO4	B	1349	5/5	0.17	3.79	51,51,53,55	5
4	SO4	C	1343	5/5	0.15	3.47	43,46,48,49	5
4	SO4	E	1343	5/5	0.12	3.35	66,68,70,71	5
4	SO4	D	1350	5/5	0.11	3.14	45,45,47,48	5
3	NO2	F	1339	3/3	0.23	2.79	69,69,69,71	0
5	ACT	E	1339	4/4	0.11	2.74	45,46,47,47	4
4	SO4	D	1337	5/5	0.13	2.66	67,68,70,72	5
4	SO4	D	1343	5/5	0.17	2.56	48,50,51,52	5
4	SO4	A	1345	5/5	0.15	2.42	55,56,57,58	5
4	SO4	A	1340	5/5	0.14	2.29	42,42,46,47	5
5	ACT	F	1336	4/4	0.12	2.16	43,44,45,45	0
4	SO4	B	1347	5/5	0.24	1.90	65,65,67,68	5
4	SO4	E	1348	5/5	0.15	1.88	42,44,46,46	5
5	ACT	D	1339	4/4	0.15	1.70	79,80,80,80	0
4	SO4	B	1348	5/5	0.22	1.67	50,54,56,57	5
4	SO4	D	1349	5/5	0.16	1.47	56,59,60,61	5
4	SO4	F	1340	5/5	0.10	1.37	52,52,53,54	5
4	SO4	C	1342	5/5	0.14	1.20	50,50,51,52	5
4	SO4	F	1345	5/5	0.18	0.99	48,50,52,52	5
4	SO4	F	1346	5/5	0.14	0.94	59,60,63,64	5
4	SO4	F	1342	5/5	0.16	0.87	56,57,59,59	5
4	SO4	A	1347	5/5	0.15	0.80	51,56,57,57	5
4	SO4	A	1344	5/5	0.15	0.78	49,50,52,54	5
4	SO4	E	1345	5/5	0.24	0.76	67,69,71,71	5
5	ACT	C	1337	4/4	0.12	0.70	41,42,43,45	0
4	SO4	E	1342	5/5	0.13	0.69	57,58,59,62	4
5	ACT	D	1338	4/4	0.10	0.60	49,49,50,51	0
4	SO4	C	1345	5/5	0.13	0.12	43,48,48,49	5
5	ACT	E	1338	4/4	0.11	0.10	52,53,53,53	0
4	SO4	D	1348	5/5	0.12	-0.16	42,44,48,50	5
5	ACT	A	1341	4/4	0.09	-0.18	39,40,41,41	0
4	SO4	E	1347	5/5	0.12	-0.25	56,58,59,59	5
4	SO4	C	1346	5/5	0.11	-0.33	53,56,57,57	5
5	ACT	B	1342	4/4	0.08	-0.33	41,42,43,43	0
4	SO4	B	1340	5/5	0.12	-0.37	48,49,52,52	5
2	CU	B	1338	1/1	0.09	-0.40	36,36,36,36	0
2	CU	E	1340	1/1	0.09	-0.49	34,34,34,34	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	NO2	B	1341	3/3	0.13	-0.56	59,59,60,60	0
4	SO4	B	1345	5/5	0.09	-0.68	49,49,50,52	5
6	NO	E	1349	2/2	0.14	-0.69	63,63,63,64	0
6	NO	C	1347	2/2	0.14	-0.75	45,45,45,46	0
3	NO2	A	1342	3/3	0.10	-0.79	69,69,69,70	0
5	ACT	B	1337	4/4	0.09	-0.85	42,42,42,43	0
2	CU	D	1341	1/1	0.09	-0.89	37,37,37,37	0
5	ACT	E	1337	4/4	0.07	-0.93	38,40,40,41	0
2	CU	F	1338	1/1	0.10	-0.96	35,35,35,35	0
2	CU	D	1340	1/1	0.08	-0.98	36,36,36,36	0
3	NO2	A	1339	3/3	0.12	-1.20	57,57,58,60	0
2	CU	F	1337	1/1	0.07	-1.33	35,35,35,35	0
2	CU	C	1338	1/1	0.08	-1.56	36,36,36,36	0
2	CU	A	1337	1/1	0.09	-2.02	35,35,35,35	0
2	CU	A	1338	1/1	0.10	-2.05	35,35,35,35	0
2	CU	B	1339	1/1	0.10	-2.22	36,36,36,36	0
2	CU	C	1339	1/1	0.09	-2.34	35,35,35,35	0
2	CU	E	1341	1/1	0.10	-2.85	36,36,36,36	0

6.5 Other polymers ⓘ

There are no such residues in this entry.