



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:37 AM UTC

PDB ID : 2BGW / pdb_00002bgw
Title : XPF from *Aeropyrum pernix*, complex with DNA
Authors : Newman, M.; Murray-Rust, J.; Lally, J.; Rudolf, J.; Fadden, A.; Knowles, P.P.; White, M.F.; McDonald, N.Q.
Deposited on : 2005-01-06
Resolution : 2.80 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

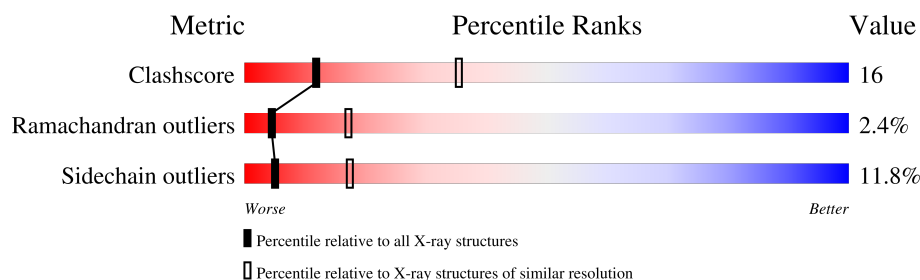
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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))
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	219	
1	B	219	
2	C	15	
3	D	15	

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
5	SO4	A	1231	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3869 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 XPF ENDONUCLEASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1677	1070	291	310	6			
1	B	200	Total	C	N	O	S	0	0	0
			1547	991	267	284	5			

- Molecule 2 is a DNA chain called 5'-D(*GP*AP*TP*CP*AP*CP*AP*GP*AP*TP*GP*CP*TP*GP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	15	Total	C	N	O	P	0	0	0
			307	147	60	86	14			

- Molecule 3 is a DNA chain called 5'-D(*TP*CP*AP*GP*CP*AP*TP*CP*TP*GP*TP*GP*AP*TP*C)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	15	Total	C	N	O	P	0	0	0
			302	146	52	90	14			

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			3	2	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	3	Total	O	0	0
			3	3		
6	B	4	Total	O	0	0
			4	4		

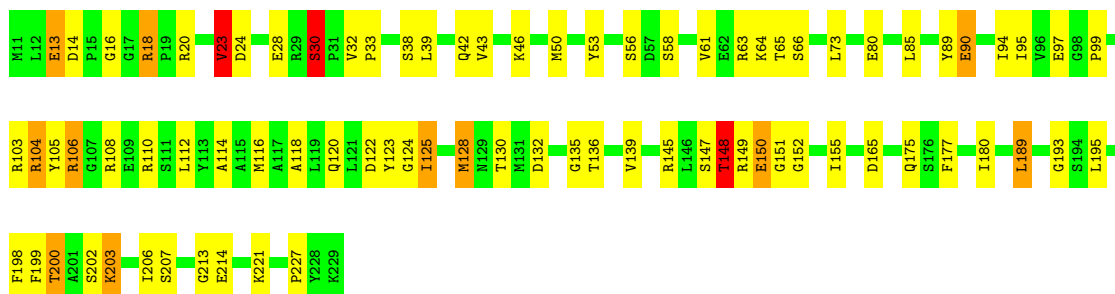
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

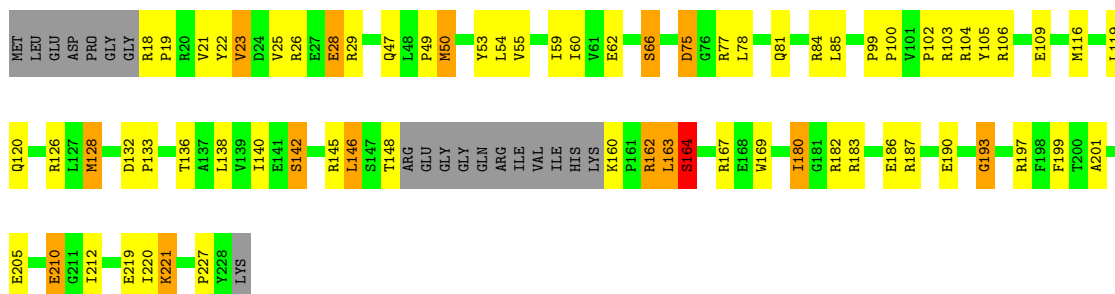
• Molecule 1: XPF ENDONUCLEASE

Chain A: 



• Molecule 1: XPF ENDONUCLEASE

Chain B: 



• Molecule 2: 5'-D(*GP*AP*TP*CP*AP*CP*AP*GP*AP*TP *GP*CP*TP*GP*A)-3'

Chain C: 



• Molecule 3: 5'-D(*TP*CP*AP*GP*CP*AP*TP*CP*TP*GP *TP*GP*AP*TP*C)-3'

Chain D: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	141.31Å 141.31Å 85.31Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.80	Depositor
% Data completeness (in resolution range)	97.3 (30.00-2.80)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.223 , 0.263	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3869	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, 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	0.91	1/1709 (0.1%)	1.14	5/2316 (0.2%)
1	B	0.92	1/1576 (0.1%)	1.15	8/2132 (0.4%)
2	C	0.66	0/345	1.80	12/531 (2.3%)
3	D	0.57	0/337	1.60	10/518 (1.9%)
All	All	0.87	2/3967 (0.1%)	1.27	35/5497 (0.6%)

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
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	89	TYR	CA-C	-5.25	1.46	1.52
1	B	163	LEU	N-CA	5.10	1.52	1.46

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	23	VAL	CB-CA-C	-10.13	97.06	110.84
2	C	2	DA	N9-C1'-C2'	8.74	126.61	113.50
2	C	2	DA	O5'-C5'-C4'	-6.70	100.75	110.80
1	B	163	LEU	N-CA-C	6.53	124.70	110.80
2	C	3	DT	C4'-C3'-O3'	-6.43	100.35	110.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	10	DT	N1-C1'-C2'	6.36	123.04	113.50
1	B	164	SER	N-CA-C	6.36	124.34	110.80
1	B	59	ILE	N-CA-C	6.07	116.35	107.37
1	A	13	GLU	N-CA-C	6.04	120.72	112.68
1	A	56	SER	N-CA-C	-6.03	100.42	108.74
1	B	162	ARG	N-CA-C	6.02	123.62	110.80
3	D	2	DC	C4'-C3'-O3'	-5.96	101.06	110.00
3	D	12	DG	P-O3'-C3'	5.91	129.07	120.20
2	C	11	DG	C2'-C3'-O3'	5.88	120.32	111.50
3	D	2	DC	P-O3'-C3'	5.82	128.93	120.20
1	B	120	GLN	N-CA-C	5.80	117.41	111.14
2	C	13	DT	N1-C1'-C2'	5.73	122.09	113.50
3	D	6	DA	O3'-P-O5'	-5.70	95.45	104.00
3	D	10	DG	O4'-C1'-N9	-5.65	99.93	108.40
2	C	3	DT	O5'-C5'-C4'	-5.61	102.39	110.80
1	B	75	ASP	CB-CA-C	-5.54	99.07	110.38
2	C	1	DG	C4'-C3'-O3'	-5.54	101.68	110.00
2	C	6	DC	C5'-C4'-O4'	5.49	117.64	109.40
1	A	80	GLU	N-CA-C	-5.47	105.23	111.14
1	A	148	THR	N-CA-C	-5.47	99.14	110.80
3	D	12	DG	C4'-C3'-O3'	5.44	118.16	110.00
3	D	5	DC	C1'-O4'-C4'	-5.37	101.65	109.70
3	D	1	DT	C5'-C4'-C3'	-5.28	106.98	114.90
2	C	2	DA	C1'-O4'-C4'	-5.24	101.83	109.70
1	B	180	ILE	N-CA-C	5.20	115.26	107.51
2	C	10	DT	O3'-P-O5'	-5.16	96.26	104.00
3	D	8	DC	N1-C1'-C2'	5.15	121.23	113.50
2	C	11	DG	P-O3'-C3'	5.11	127.86	120.20
1	B	212	ILE	CB-CA-C	-5.09	104.60	110.41
3	D	9	DT	P-O3'-C3'	5.04	127.75	120.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	150	GLU	Peptide
1	B	160	LYS	Peptide

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	1677	0	1656	59	0
1	B	1547	0	1540	52	0
2	C	307	0	170	11	0
3	D	302	0	172	5	0
4	A	1	0	0	0	0
5	A	13	0	0	0	3
5	B	10	0	0	1	0
5	D	5	0	0	1	0
6	A	3	0	0	0	0
6	B	4	0	0	0	0
All	All	3869	0	3538	115	3

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 16.

All (115) 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:187:ARG:HD3	1:B:210:GLU:HG2	1.53	0.90
1:B:187:ARG:CD	1:B:210:GLU:HG2	2.01	0.90
1:B:66:SER:OG	1:B:102:PRO:HG3	1.80	0.82
1:A:148:THR:O	1:A:149:ARG:HB2	1.82	0.79
1:B:50:MET:HG2	1:B:84:ARG:NH2	1.97	0.79
1:A:50:MET:HE2	1:A:85:LEU:HA	1.63	0.78
1:A:95:ILE:HG21	1:A:136:THR:HG23	1.66	0.77
1:A:14:ASP:O	1:A:149:ARG:HA	1.85	0.75
1:B:201:ALA:O	1:B:221:LYS:HE3	1.87	0.74
1:A:116:MET:HB3	1:B:128:MET:HE1	1.67	0.74
1:B:50:MET:HE3	1:B:85:LEU:HD23	1.70	0.74
1:B:187:ARG:HD3	1:B:210:GLU:CG	2.18	0.73
1:B:180:ILE:HD11	1:B:220:ILE:HD11	1.70	0.72
1:A:128:MET:HE1	1:B:116:MET:HB3	1.74	0.69
1:A:16:GLY:HA3	1:A:148:THR:CG2	2.22	0.69
1:A:58:SER:HB2	1:A:90:GLU:HG2	1.73	0.69
1:B:187:ARG:HD2	1:B:210:GLU:HG2	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:TYR:HB2	1:A:125:ILE:HD12	1.75	0.68
1:A:18:ARG:HG2	1:A:18:ARG:HH11	1.60	0.67
1:A:13:GLU:H	1:A:151:GLY:HA2	1.59	0.66
1:B:132:ASP:HB2	1:B:133:PRO:CD	2.26	0.66
1:B:210:GLU:H	1:B:210:GLU:CD	2.06	0.63
1:A:128:MET:HE1	1:B:116:MET:CB	2.29	0.62
1:A:94:ILE:HD12	1:A:125:ILE:HG21	1.81	0.62
1:B:103:ARG:HA	1:B:106:ARG:HE	1.64	0.61
1:A:18:ARG:HG2	1:A:18:ARG:NH1	2.15	0.60
2:C:12:DC:H2'	2:C:13:DT:C6	2.37	0.60
1:A:147:SER:C	1:A:148:THR:O	2.40	0.60
1:A:114:ALA:HA	1:B:142:SER:OG	2.01	0.59
1:A:165:ASP:N	1:A:165:ASP:OD1	2.32	0.59
1:B:182:ARG:O	1:B:186:GLU:HG2	2.03	0.58
1:A:63:ARG:NH2	1:A:97:GLU:OE2	2.35	0.58
1:A:199:PHE:CD2	1:B:199:PHE:CD2	2.92	0.58
1:B:164:SER:HA	1:B:169:TRP:HE1	1.70	0.57
1:A:103:ARG:O	1:A:106:ARG:HD2	2.04	0.57
1:A:124:GLY:HA3	3:D:1:DT:C6	2.41	0.56
1:A:193:GLY:O	1:B:227:PRO:HA	2.05	0.56
1:A:103:ARG:O	1:A:105:TYR:N	2.38	0.56
1:A:13:GLU:HG3	1:A:152:GLY:HA3	1.89	0.55
1:A:124:GLY:O	1:B:126:ARG:NH2	2.40	0.55
1:A:148:THR:HG22	1:A:149:ARG:N	2.22	0.55
1:B:49:PRO:O	1:B:50:MET:HG3	2.07	0.54
1:A:148:THR:HG22	1:A:149:ARG:H	1.71	0.54
1:B:145:ARG:O	1:B:148:THR:HB	2.08	0.53
2:C:4:DC:H2''	2:C:5:DA:C8	2.43	0.53
1:B:62:GLU:CD	1:B:81:GLN:NE2	2.67	0.52
1:B:23:VAL:HG22	1:B:53:TYR:CD2	2.44	0.52
1:B:142:SER:O	1:B:146:LEU:HB2	2.09	0.52
1:A:148:THR:O	1:A:149:ARG:CB	2.49	0.52
1:A:104:ARG:O	1:A:104:ARG:HG3	2.10	0.52
2:C:7:DA:H2''	2:C:8:DG:OP2	2.10	0.51
1:B:183:ARG:CB	1:B:187:ARG:HH21	2.22	0.51
1:A:23:VAL:HG13	1:A:53:TYR:CE2	2.46	0.51
2:C:3:DT:H2''	2:C:4:DC:O5'	2.11	0.51
2:C:2:DA:H2'	2:C:3:DT:C6	2.45	0.51
1:B:136:THR:O	1:B:140:ILE:HG13	2.11	0.50
1:A:120:GLN:HE22	1:B:126:ARG:HB3	1.76	0.50
2:C:12:DC:H2''	2:C:13:DT:H5'	1.91	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:SER:O	1:A:203:LYS:C	2.54	0.50
1:A:30:SER:HB3	1:A:32:VAL:H	1.77	0.50
2:C:12:DC:H2''	2:C:13:DT:C5'	2.41	0.50
1:B:50:MET:CE	1:B:85:LEU:HD23	2.40	0.50
1:B:103:ARG:HA	1:B:106:ARG:NE	2.27	0.50
1:B:103:ARG:HB2	1:B:106:ARG:HE	1.76	0.49
1:B:50:MET:CG	1:B:84:ARG:NH2	2.74	0.49
1:B:132:ASP:HB2	1:B:133:PRO:HD2	1.94	0.48
1:A:32:VAL:HB	1:A:33:PRO:HD3	1.94	0.48
2:C:2:DA:H2'	2:C:3:DT:H6	1.79	0.47
3:D:8:DC:H1'	3:D:9:DT:H5'	1.97	0.47
1:A:103:ARG:C	1:A:105:TYR:H	2.23	0.47
1:A:104:ARG:HG2	1:A:105:TYR:CD1	2.50	0.47
1:A:130:THR:HG21	1:A:136:THR:HA	1.97	0.47
1:A:65:THR:O	1:A:66:SER:C	2.58	0.46
1:A:53:TYR:HB2	1:A:61:VAL:HB	1.96	0.46
1:A:104:ARG:O	1:A:104:ARG:CG	2.62	0.46
1:A:227:PRO:HA	1:B:193:GLY:O	2.15	0.46
1:A:20:ARG:HG3	1:A:42:GLN:OE1	2.16	0.46
1:A:66:SER:CB	1:A:99:PRO:O	2.64	0.45
1:B:22:TYR:HB2	1:B:54:LEU:HB3	1.98	0.45
1:B:62:GLU:CD	1:B:81:GLN:HE22	2.25	0.45
1:B:49:PRO:C	1:B:50:MET:HG3	2.42	0.44
1:A:24:ASP:HA	1:A:46:LYS:O	2.18	0.44
1:B:99:PRO:HA	1:B:100:PRO:HD3	1.92	0.44
1:A:16:GLY:HA3	1:A:148:THR:HG22	1.98	0.44
1:B:78:LEU:HD21	1:B:119:LEU:HD11	1.98	0.44
1:A:118:ALA:O	1:A:122:ASP:HB2	2.18	0.43
1:B:23:VAL:HG22	1:B:53:TYR:CE2	2.53	0.43
1:A:189:LEU:HA	1:A:189:LEU:HD23	1.53	0.43
1:A:198:PHE:C	1:A:200:THR:H	2.26	0.43
1:A:110:ARG:HD2	1:B:138:LEU:HD22	2.00	0.43
1:B:197:ARG:NH2	1:B:205:GLU:OE2	2.50	0.43
1:B:18:ARG:HA	1:B:19:PRO:HD2	1.68	0.42
1:A:177:PHE:HB2	1:A:180:ILE:HD12	2.01	0.42
1:B:103:ARG:C	1:B:105:TYR:H	2.26	0.42
1:B:50:MET:HB3	5:B:1229:SO4:O1	2.19	0.42
1:A:120:GLN:OE1	1:B:126:ARG:NE	2.44	0.42
1:A:135:GLY:O	1:A:139:VAL:HG23	2.20	0.42
2:C:10:DT:H2''	2:C:11:DG:C8	2.53	0.42
2:C:2:DA:H2''	2:C:3:DT:O5'	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:ASP:HB2	1:B:77:ARG:H	1.84	0.41
1:B:180:ILE:CD1	1:B:220:ILE:HD11	2.44	0.41
1:A:151:GLY:HA3	1:A:155:ILE:HG21	2.02	0.41
1:A:73:LEU:HD23	1:A:112:LEU:HD23	2.02	0.41
1:A:112:LEU:HD23	1:A:112:LEU:HA	1.94	0.41
1:B:197:ARG:HH22	1:B:205:GLU:CD	2.29	0.41
2:C:11:DG:H1'	2:C:12:DC:C6	2.55	0.41
1:B:25:VAL:O	1:B:28:GLU:HB2	2.20	0.41
3:D:9:DT:H2''	3:D:10:DG:C8	2.56	0.41
1:A:213:GLY:O	1:A:214:GLU:C	2.64	0.41
3:D:1:DT:H5''	5:D:1016:SO4:O1	2.21	0.41
1:A:23:VAL:HG22	1:A:43:VAL:HG13	2.02	0.40
3:D:10:DG:OP2	3:D:10:DG:H8	2.04	0.40
1:A:18:ARG:HH11	1:A:18:ARG:CG	2.29	0.40
1:A:206:ILE:O	1:A:206:ILE:HG22	2.22	0.40
1:B:103:ARG:CA	1:B:106:ARG:HE	2.32	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:1231:SO4:S	5:A:1231:SO4:O1[4_556]	1.44	0.76
5:A:1231:SO4:S	5:A:1231:SO4:O2[4_556]	1.47	0.73
5:A:1231:SO4:O1	5:A:1231:SO4:O1[4_556]	2.03	0.17

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	217/219 (99%)	200 (92%)	12 (6%)	5 (2%)	5	18
1	B	196/219 (90%)	185 (94%)	6 (3%)	5 (3%)	4	15

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	413/438 (94%)	385 (93%)	18 (4%)	10 (2%)	4	17

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	104	ARG
1	B	163	LEU
1	B	164	SER
1	A	28	GLU
1	B	162	ARG
1	A	148	THR
1	B	104	ARG
1	B	193	GLY
1	A	203	LYS
1	A	30	SER

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	172/190 (90%)	152 (88%)	20 (12%)	5	18
1	B	159/190 (84%)	140 (88%)	19 (12%)	5	17
All	All	331/380 (87%)	292 (88%)	39 (12%)	5	17

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ARG
1	A	23	VAL
1	A	30	SER
1	A	38	SER
1	A	39	LEU
1	A	64	LYS
1	A	90	GLU
1	A	106	ARG

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Mol	Chain	Res	Type
1	A	108	ARG
1	A	125	ILE
1	A	128	MET
1	A	132	ASP
1	A	145	ARG
1	A	150	GLU
1	A	175	GLN
1	A	189	LEU
1	A	195	LEU
1	A	200	THR
1	A	207	SER
1	A	221	LYS
1	B	21	VAL
1	B	23	VAL
1	B	26	ARG
1	B	28	GLU
1	B	29	ARG
1	B	47	GLN
1	B	50	MET
1	B	55	VAL
1	B	60	ILE
1	B	66	SER
1	B	109	GLU
1	B	128	MET
1	B	142	SER
1	B	146	LEU
1	B	167	ARG
1	B	190	GLU
1	B	210	GLU
1	B	219	GLU
1	B	221	LYS

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

Mol	Chain	Res	Type
1	A	129	ASN
1	B	42	GLN
1	B	88	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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	SO4	D	1016	-	4,4,4	0.25	0	6,6,6	0.41	0
5	SO4	A	1233	-	4,4,4	0.39	0	6,6,6	0.65	0
5	SO4	A	1232	-	4,4,4	0.26	0	6,6,6	0.42	0
5	SO4	B	1229	-	4,4,4	0.27	0	6,6,6	0.42	0
5	SO4	B	1230	-	4,4,4	0.40	0	6,6,6	0.57	0
5	SO4	A	1231	-	2,2,4	0.77	0	1,1,6	0.68	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.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	1016	SO4	1	0
5	B	1229	SO4	1	0
5	A	1231	SO4	0	3

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.