Full wwPDB X-ray Structure Validation Report (i)

Apr 29, 2024 – 06:07 am BST

PDB ID	:	2VTW
Title	:	Structure of the C-terminal head domain of the fowl adenovirus type 1 short fibre
Authors	:	ElBakkouri, M.; Seiradake, E.; Cusack, S.; Ruigrok, R.W.H.; Schoehn, G.
Deposited on	:	2008-05-16
Resolution	:	2.00 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:

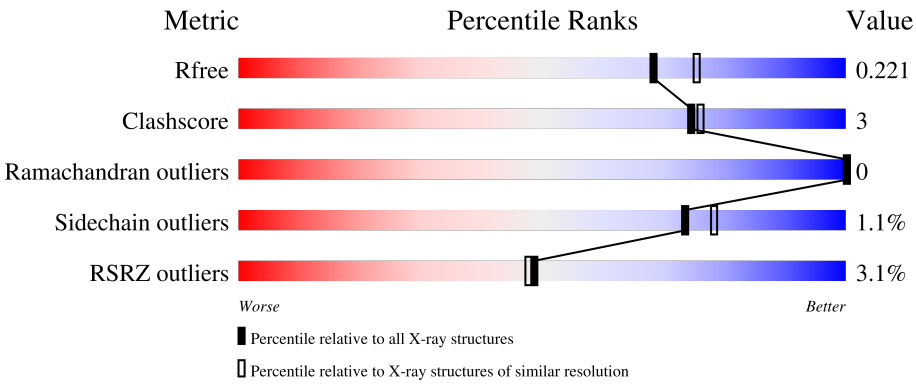
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	205	6% 92% 7% •
1	B	205	% 94% 5% •
1	C	205	8% 90% 9% •
1	D	205	% 92% 8%
1	E	205	% 93% 6%

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Mol	Chain	Length	Quality of chain
1	F	205	% 96% .

2 Entry composition (i)

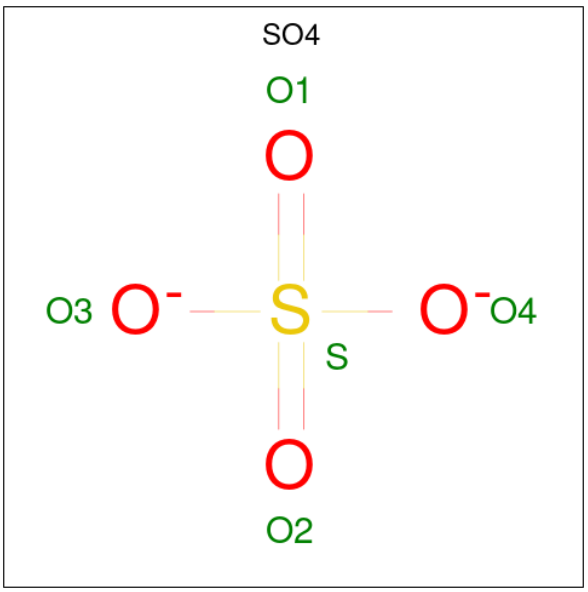
There are 4 unique types of molecules in this entry. The entry contains 10479 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 FIBER PROTEIN 2.

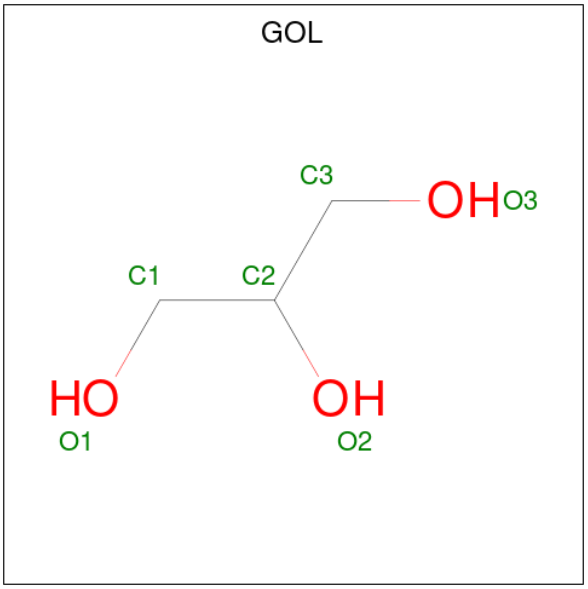
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	205	1541	972	258	302	9	0	1	0
1	B	205	1540	973	257	301	9	0	1	0
1	C	205	1536	969	257	301	9	0	0	0
1	D	205	1545	976	258	302	9	0	2	0
1	E	205	1542	973	257	303	9	0	2	0
1	F	205	1536	969	257	301	9	0	0	0

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	B	1	6	3	3	0	0
3	E	1	6	3	3	0	0

- Molecule 4 is water.

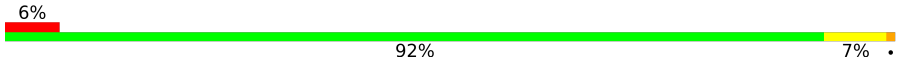
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	139	139	139	0	0
			Total	O		
4	B	253	253	253	0	0
			Total	O		
4	C	107	107	107	0	0
			Total	O		
4	D	208	208	208	0	0
			Total	O		
4	E	256	256	256	0	0
			Total	O		
4	F	214	214	214	0	0

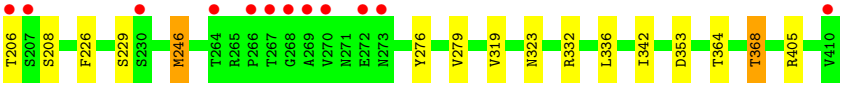
3 Residue-property plots (i)

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: FIBER PROTEIN 2

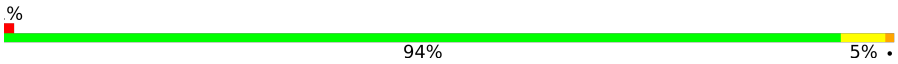
Chain A:





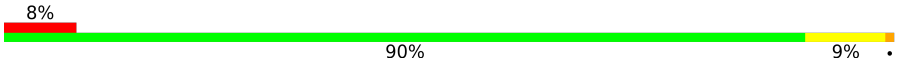
- Molecule 1: FIBER PROTEIN 2

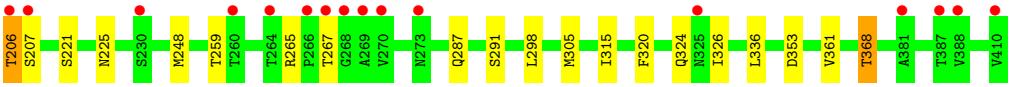
Chain B:



- Molecule 1: FIBER PROTEIN 2

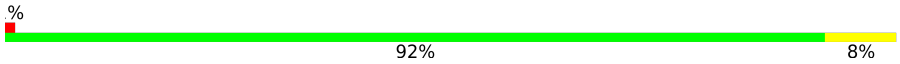
Chain C:

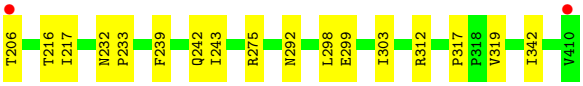




- Molecule 1: FIBER PROTEIN 2

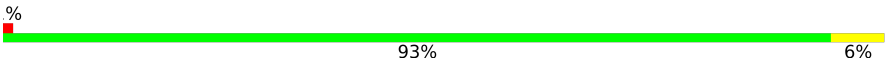
Chain D:

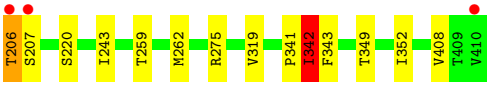




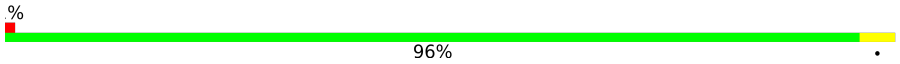
- Molecule 1: FIBER PROTEIN 2

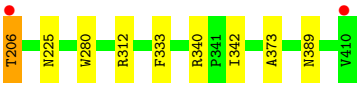
Chain E:





- Molecule 1: FIBER PROTEIN 2

Chain F: 



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	235.75Å 235.75Å 61.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.57 – 2.00 47.56 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.57-2.00) 99.6 (47.56-2.00)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.31 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.175 , 0.214 0.183 , 0.221	Depositor DCC
R_{free} test set	5865 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	19.0	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 49.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	10479	wwPDB-VP
Average B, all atoms (Å^2)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.07 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.4307e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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: GOL, 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	0.70	1/1582 (0.1%)	0.74	3/2169 (0.1%)
1	B	0.83	2/1581 (0.1%)	0.76	1/2168 (0.0%)
1	C	0.67	1/1574 (0.1%)	0.69	0/2158
1	D	0.79	1/1589 (0.1%)	0.76	3/2179 (0.1%)
1	E	0.77	1/1586 (0.1%)	0.76	2/2175 (0.1%)
1	F	0.82	1/1574 (0.1%)	0.74	1/2158 (0.0%)
All	All	0.77	7/9486 (0.1%)	0.74	10/13007 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	206	THR	C-N	-15.41	0.98	1.34
1	D	206	THR	C-N	-13.24	1.03	1.34
1	B	206	THR	C-N	-12.30	1.05	1.34
1	A	206	THR	C-N	-8.85	1.13	1.34
1	C	206	THR	C-N	-7.89	1.15	1.34
1	B	220	SER	CB-OG	5.37	1.49	1.42
1	E	206	THR	C-N	-5.26	1.22	1.34

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	206	THR	O-C-N	6.99	133.89	122.70
1	B	275	ARG	NE-CZ-NH1	6.55	123.58	120.30
1	A	206	THR	CA-C-N	-6.32	103.29	117.20
1	E	275	ARG	NE-CZ-NH1	5.61	123.11	120.30
1	E	342	ILE	CG1-CB-CG2	-5.32	99.70	111.40
1	F	312	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	D	312	ARG	NE-CZ-NH1	5.25	122.93	120.30
1	D	275	ARG	NE-CZ-NH1	5.09	122.85	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	405	ARG	NE-CZ-NH1	5.05	122.83	120.30
1	D	312	ARG	NE-CZ-NH2	-5.03	117.78	120.30

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	1541	0	1501	9	0
1	B	1540	0	1504	7	0
1	C	1536	0	1495	14	0
1	D	1545	0	1510	15	0
1	E	1542	0	1506	13	0
1	F	1536	0	1495	7	0
2	A	5	0	0	0	0
2	B	10	0	0	0	0
2	C	5	0	0	1	0
2	D	10	0	0	0	0
2	E	10	0	0	0	0
2	F	10	0	0	0	0
3	B	6	0	8	0	0
3	E	6	0	7	0	0
4	A	139	0	0	3	0
4	B	253	0	0	1	0
4	C	107	0	0	3	0
4	D	208	0	0	5	0
4	E	256	0	0	2	0
4	F	214	0	0	5	0
All	All	10479	0	9026	58	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 3.

All (58) 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:C:368:THR:HG23	4:C:2091:HOH:O	1.60	0.98
1:D:292[B]:ASN:ND2	4:D:2072:HOH:O	1.83	0.97
1:F:206:THR:HB	4:F:2001:HOH:O	1.69	0.92
1:B:286:THR:HG23	4:B:2248:HOH:O	1.82	0.77
1:E:206:THR:HG23	1:E:207:SER:N	2.01	0.75
1:E:206:THR:HG23	1:E:207:SER:H	1.53	0.73
1:F:206:THR:CB	4:F:2001:HOH:O	2.31	0.73
1:A:368:THR:HG23	4:A:2118:HOH:O	1.89	0.73
1:C:315:ILE:HD12	1:D:342:ILE:HG13	1.77	0.67
1:D:242:GLN:C	1:D:243:ILE:HD12	2.16	0.67
1:A:319:VAL:HG12	4:A:2074:HOH:O	1.96	0.65
1:D:217:ILE:HD13	1:D:239:PHE:CE2	2.33	0.64
1:D:319[B]:VAL:HG13	4:D:2116:HOH:O	1.98	0.62
1:D:319[B]:VAL:HG12	4:D:2117:HOH:O	1.99	0.61
1:E:206:THR:CG2	1:E:207:SER:H	2.13	0.61
1:C:368:THR:CG2	4:C:2091:HOH:O	2.30	0.61
1:A:368:THR:CG2	4:A:2118:HOH:O	2.45	0.61
1:F:389:ASN:ND2	4:F:2193:HOH:O	2.18	0.59
1:E:206:THR:CG2	1:E:207:SER:N	2.66	0.58
1:F:225:ASN:HB3	4:F:2018:HOH:O	2.03	0.58
1:C:248:MET:CE	1:D:243:ILE:HD13	2.35	0.56
1:A:323:ASN:OD1	1:A:364:THR:HA	2.07	0.53
1:A:226:PHE:CD1	1:A:279:VAL:HG22	2.44	0.52
1:F:280:TRP:CZ3	1:F:373:ALA:HB2	2.45	0.52
1:E:342:ILE:HG12	1:F:333:PHE:CD1	2.47	0.50
1:C:259:THR:HB	1:C:265:ARG:NH2	2.28	0.49
1:C:320:PHE:HB3	1:C:361:VAL:HG22	1.95	0.49
1:E:408:VAL:CG1	4:E:2247:HOH:O	2.60	0.49
1:D:216:THR:OG1	4:D:2007:HOH:O	2.17	0.48
1:C:248:MET:HE3	1:D:243:ILE:HD13	1.96	0.48
1:B:323:ASN:OD1	1:B:364:THR:HA	2.13	0.47
1:C:248:MET:HE2	1:D:243:ILE:HD13	1.96	0.47
1:C:287:GLN:HG3	2:C:1411:SO4:O1	2.13	0.47
1:E:341:PRO:HG3	1:E:352:ILE:HD12	1.97	0.47
1:E:319:VAL:HG13	4:E:2138:HOH:O	2.14	0.46
1:A:208:SER:HB2	1:A:246:MET:HG3	1.97	0.46
1:C:324:GLN:NE2	4:C:2059:HOH:O	2.48	0.45
1:D:243:ILE:HD12	1:D:243:ILE:N	2.32	0.44
1:C:206:THR:HG22	1:C:207:SER:N	2.31	0.44
1:B:243:ILE:HD11	4:D:2080:HOH:O	2.16	0.43
1:E:243:ILE:HD11	4:F:2077:HOH:O	2.18	0.43
1:B:343:PHE:HE1	1:B:349:THR:HG1	1.64	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:342:ILE:HD12	1:E:342:ILE:HG21	1.74	0.42
1:B:213:THR:O	1:B:296:GLY:HA3	2.20	0.42
1:E:343:PHE:HE1	1:E:349:THR:HG1	1.65	0.42
1:A:336:LEU:HD11	1:A:353:ASP:HB3	2.01	0.42
1:A:229:SER:HB3	1:A:276:TYR:HB2	2.02	0.42
1:A:332:ARG:O	1:F:340:ARG:HD2	2.20	0.42
1:C:336:LEU:HD11	1:C:353:ASP:HB3	2.01	0.42
1:B:342:ILE:HD11	1:D:317:PRO:HA	2.02	0.41
1:E:259:THR:HA	1:E:262:MET:SD	2.60	0.41
1:D:298:LEU:HB3	1:D:303:ILE:HG22	2.02	0.41
1:E:341:PRO:HG3	1:E:352:ILE:CD1	2.50	0.41
1:B:206:THR:O	1:B:207:SER:C	2.59	0.41
1:D:232:ASN:HA	1:D:233:PRO:HD3	1.96	0.40
1:D:242:GLN:O	1:D:243:ILE:HD12	2.20	0.40
1:C:225:ASN:O	1:C:291:SER:HA	2.22	0.40
1:C:298:LEU:HD21	1:C:305:MET:SD	2.60	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	204/205 (100%)	198 (97%)	6 (3%)	0	100	100
1	B	204/205 (100%)	199 (98%)	5 (2%)	0	100	100
1	C	203/205 (99%)	198 (98%)	5 (2%)	0	100	100
1	D	205/205 (100%)	198 (97%)	7 (3%)	0	100	100
1	E	205/205 (100%)	201 (98%)	4 (2%)	0	100	100
1	F	203/205 (99%)	198 (98%)	5 (2%)	0	100	100
All	All	1224/1230 (100%)	1192 (97%)	32 (3%)	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	177/176 (101%)	174 (98%)	3 (2%)	60	65
1	B	177/176 (101%)	176 (99%)	1 (1%)	86	90
1	C	176/176 (100%)	172 (98%)	4 (2%)	50	53
1	D	178/176 (101%)	177 (99%)	1 (1%)	86	90
1	E	178/176 (101%)	175 (98%)	3 (2%)	60	65
1	F	176/176 (100%)	175 (99%)	1 (1%)	86	90
All	All	1062/1056 (101%)	1049 (99%)	13 (1%)	73	76

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	246	MET
1	A	342	ILE
1	A	368	THR
1	B	220	SER
1	C	221	SER
1	C	267	THR
1	C	326	ILE
1	C	368	THR
1	D	299	GLU
1	E	220[A]	SER
1	E	220[B]	SER
1	E	342	ILE
1	F	342	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	235	ASN

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Mol	Chain	Res	Type
1	E	366	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

12 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	$\#\|Z\| > 2$
2	SO4	A	1411	-	4,4,4	0.24	0	6,6,6	0.59	0
2	SO4	F	1412	-	4,4,4	0.15	0	6,6,6	0.14	0
2	SO4	C	1411	-	4,4,4	0.25	0	6,6,6	0.43	0
2	SO4	D	1412	-	4,4,4	0.18	0	6,6,6	0.17	0
2	SO4	F	1411	-	4,4,4	0.21	0	6,6,6	1.03	0
2	SO4	D	1411	-	4,4,4	0.31	0	6,6,6	0.74	0
2	SO4	B	1411	-	4,4,4	0.22	0	6,6,6	0.42	0
3	GOL	E	1413	-	5,5,5	0.99	0	5,5,5	0.70	0
2	SO4	E	1412	-	4,4,4	0.23	0	6,6,6	0.34	0
2	SO4	E	1411	-	4,4,4	0.24	0	6,6,6	0.46	0
3	GOL	B	1413	-	5,5,5	0.53	0	5,5,5	0.48	0
2	SO4	B	1412	-	4,4,4	0.14	0	6,6,6	0.47	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	GOL	B	1413	-	-	2/4/4/4	-
3	GOL	E	1413	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	1413	GOL	C1-C2-C3-O3
3	B	1413	GOL	O1-C1-C2-O2
3	E	1413	GOL	O1-C1-C2-O2
3	B	1413	GOL	O1-C1-C2-C3
3	E	1413	GOL	O1-C1-C2-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1411	SO4	1	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	1
1	A	1
1	B	1
1	D	1
1	F	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	206:THR	C	207:SER	N	1.15
1	A	206:THR	C	207:SER	N	1.13
1	B	206:THR	C	207:SER	N	1.05
1	D	206:THR	C	207:SER	N	1.03
1	F	206:THR	C	207:SER	N	0.98

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	205/205 (100%)	0.06	12 (5%)	22	21	14, 26, 44, 55	0
1	B	205/205 (100%)	-0.44	3 (1%)	73	72	10, 15, 26, 36	0
1	C	205/205 (100%)	0.22	16 (7%)	13	12	17, 28, 47, 59	0
1	D	205/205 (100%)	-0.46	2 (0%)	82	81	10, 16, 28, 36	0
1	E	205/205 (100%)	-0.42	3 (1%)	73	72	9, 14, 25, 37	0
1	F	205/205 (100%)	-0.51	2 (0%)	82	81	8, 15, 27, 35	0
All	All	1230/1230 (100%)	-0.26	38 (3%)	49	48	8, 18, 38, 59	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	206	THR	7.1
1	B	206	THR	6.9
1	E	410	VAL	5.8
1	B	410	VAL	5.5
1	C	264	THR	5.4
1	C	270	VAL	5.2
1	D	206	THR	5.2
1	C	269	ALA	5.0
1	A	206	THR	4.9
1	C	206	THR	4.3
1	F	206	THR	4.1
1	D	410	VAL	4.0
1	C	410	VAL	4.0
1	F	410	VAL	3.7
1	A	264	THR	3.7
1	A	267	THR	3.3
1	B	207	SER	3.1
1	A	270	VAL	2.9
1	C	207	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	268	GLY	2.8
1	C	267	THR	2.7
1	A	266	PRO	2.7
1	A	230	SER	2.7
1	C	266	PRO	2.6
1	C	387	THR	2.5
1	C	273	ASN	2.4
1	A	269	ALA	2.3
1	A	273	ASN	2.2
1	C	268	GLY	2.2
1	A	410	VAL	2.2
1	E	207	SER	2.1
1	A	272	GLU	2.1
1	C	388	VAL	2.1
1	C	381	ALA	2.1
1	C	325	ASN	2.0
1	C	260	THR	2.0
1	A	207	SER	2.0
1	C	230	SER	2.0

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

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

6.3 Carbohydrates (i)

There are no 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
2	SO4	D	1412	5/5	0.78	0.20	74,75,76,76	0
2	SO4	F	1411	5/5	0.90	0.16	38,41,42,44	0
2	SO4	C	1411	5/5	0.92	0.15	45,49,51,52	0
3	GOL	E	1413	6/6	0.94	0.12	17,23,24,26	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	SO4	F	1412	5/5	0.95	0.17	61,62,64,64	0
2	SO4	B	1412	5/5	0.96	0.18	48,48,50,50	0
2	SO4	D	1411	5/5	0.97	0.09	22,27,28,30	0
2	SO4	A	1411	5/5	0.97	0.10	34,36,39,40	0
2	SO4	E	1411	5/5	0.97	0.16	40,42,44,44	0
3	GOL	B	1413	6/6	0.98	0.08	17,21,23,25	0
2	SO4	B	1411	5/5	0.98	0.17	33,34,36,40	0
2	SO4	E	1412	5/5	0.99	0.12	30,32,34,35	0

6.5 Other polymers (i)

There are no such residues in this entry.