Full wwPDB X-ray Structure Validation Report (i)

Nov 1, 2023 – 01:47 AM EDT

PDB ID	:	3T9K
Title	:	Crystal Structure of ACAP1 C-portion mutant S554D fused with integrin beta1 peptide
Authors	:	Sun, F.; Pang, X.; Zhang, K.; Ma, J.; Zhou, Q.
Deposited on	:	2011-08-03
Resolution	:	2.30 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:

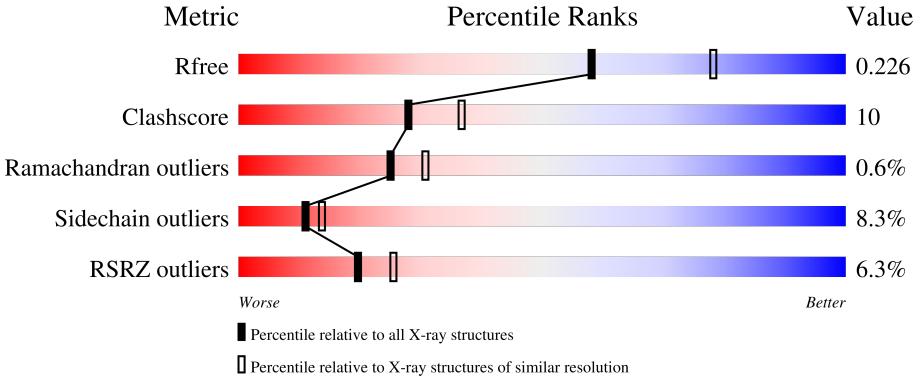
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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	390	5% 57% 13% • 28%
1	B	390	4% 58% 10% • 30%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4301 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 Arf-GAP with coiled-coil, ANK repeat and PH domain-containing protein 1,Peptide from Integrin beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	281	2106	1314	382	396	14	0	0	0
1	B	273	2008	1249	367	378	14	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	373	GLY	-	expression tag	UNP Q15027
A	374	PRO	-	expression tag	UNP Q15027
A	375	LEU	-	expression tag	UNP Q15027
A	376	GLY	-	expression tag	UNP Q15027
A	377	SER	-	expression tag	UNP Q15027
A	554	ASP	SER	engineered mutation	UNP Q15027
A	741	GLY	-	linker	UNP P05556
A	742	SER	-	linker	UNP P05556
A	743	SER	-	linker	UNP P05556
A	744	ASN	-	linker	UNP P05556
A	745	SER	-	linker	UNP P05556
A	746	GLY	-	linker	UNP P05556
A	747	SER	-	linker	UNP P05556
A	748	ASN	-	linker	UNP P05556
A	749	SER	-	linker	UNP P05556
A	750	GLY	-	linker	UNP P05556
B	373	GLY	-	expression tag	UNP Q15027
B	374	PRO	-	expression tag	UNP Q15027
B	375	LEU	-	expression tag	UNP Q15027
B	376	GLY	-	expression tag	UNP Q15027
B	377	SER	-	expression tag	UNP Q15027
B	554	ASP	SER	engineered mutation	UNP Q15027
B	741	GLY	-	linker	UNP P05556
B	742	SER	-	linker	UNP P05556

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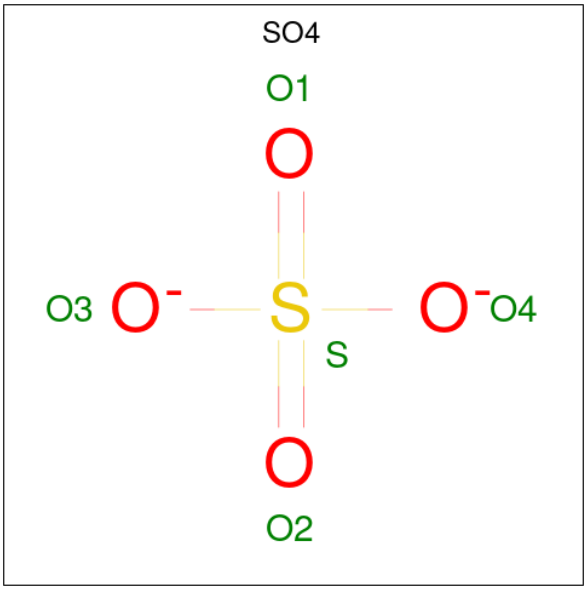
Chain	Residue	Modelled	Actual	Comment	Reference
B	743	SER	-	linker	UNP P05556
B	744	ASN	-	linker	UNP P05556
B	745	SER	-	linker	UNP P05556
B	746	GLY	-	linker	UNP P05556
B	747	SER	-	linker	UNP P05556
B	748	ASN	-	linker	UNP P05556
B	749	SER	-	linker	UNP P05556
B	750	GLY	-	linker	UNP P05556

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

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

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



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

- Molecule 4 is water.

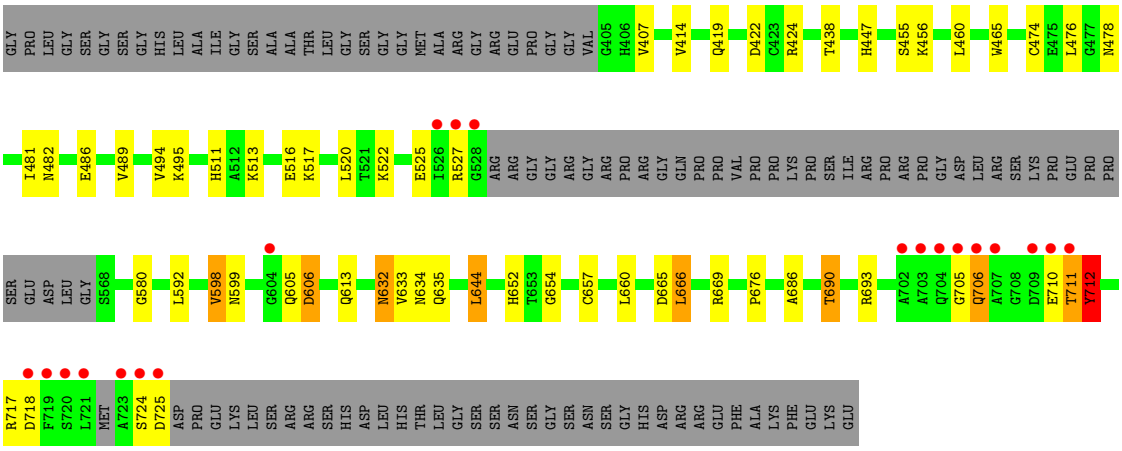
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	104	104	104	0	0
			Total	O		
4	B	71	71	71	0	0

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.

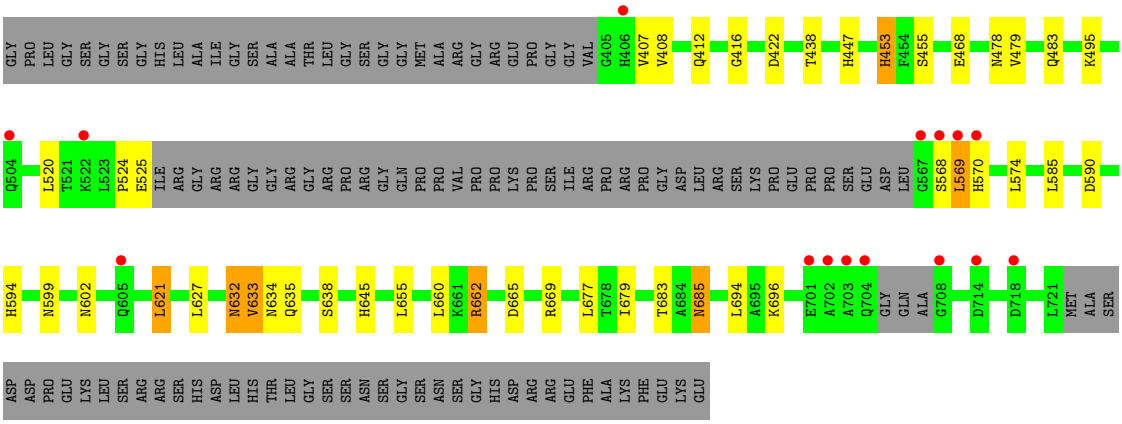
• Molecule 1: Arf-GAP with coiled-coil, ANK repeat and PH domain-containing protein 1,Peptide from Integrin beta-1

Chain A:



• Molecule 1: Arf-GAP with coiled-coil, ANK repeat and PH domain-containing protein 1,Peptide from Integrin beta-1

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	108.26Å 164.87Å 41.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.30 38.56 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.4 (40.00-2.30) 98.5 (38.56-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.63 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.189 , 0.225 0.189 , 0.226	Depositor DCC
R_{free} test set	1708 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	37.6	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 37.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4301	wwPDB-VP
Average B, all atoms (Å^2)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.40% 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, ZN

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	0.55	1/2144 (0.0%)	0.69	0/2907
1	B	0.56	0/2043	0.66	0/2773
All	All	0.56	1/4187 (0.0%)	0.68	0/5680

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	657	CYS	CB-SG	-5.44	1.73	1.81

There are no bond angle outliers.

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	2106	0	2080	50	0
1	B	2008	0	1955	31	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	104	0	0	2	0
4	B	71	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4301	0	4035	81	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 10.

All (81) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:654:GLY:H	1:A:725:ASP:CB	1.51	1.24
1:B:679:ILE:O	1:B:683:THR:HG22	1.42	1.18
1:A:654:GLY:N	1:A:725:ASP:HB2	1.59	1.18
1:A:711:THR:HA	1:A:712:TYR:CB	1.82	1.10
1:A:654:GLY:HA3	1:A:725:ASP:HB3	1.37	1.07
1:A:711:THR:CA	1:A:712:TYR:HB2	1.84	1.07
1:A:654:GLY:H	1:A:725:ASP:HB2	0.85	1.00
1:B:634:ASN:HD21	1:B:665:ASP:H	1.07	0.96
1:A:654:GLY:N	1:A:725:ASP:CB	2.22	0.92
1:A:654:GLY:CA	1:A:725:ASP:HB3	2.00	0.91
1:A:634:ASN:HD21	1:A:665:ASP:H	1.25	0.84
1:A:711:THR:HA	1:A:712:TYR:HB2	0.91	0.84
1:B:479:VAL:O	1:B:483:GLN:HG3	1.79	0.83
1:A:599:ASN:HD21	1:A:632:ASN:H	1.24	0.81
1:A:486:GLU:O	1:A:489:VAL:HG12	1.83	0.78
1:B:599:ASN:HB3	1:B:635:GLN:HE22	1.50	0.77
1:B:422:ASP:OD1	1:B:447:HIS:HE1	1.69	0.76
1:B:634:ASN:ND2	1:B:665:ASP:H	1.82	0.75
1:B:438:THR:H	1:B:478:ASN:HD21	1.34	0.74
1:B:632:ASN:ND2	1:B:634:ASN:H	1.88	0.71
1:A:654:GLY:CA	1:A:725:ASP:CB	2.66	0.71
1:B:569:LEU:CD1	1:B:570:HIS:H	2.04	0.70
1:B:453:HIS:H	1:B:453:HIS:CD2	2.09	0.70
1:A:634:ASN:ND2	1:A:665:ASP:H	1.91	0.69
1:B:599:ASN:HD21	1:B:632:ASN:H	1.40	0.69
1:A:438:THR:H	1:A:478:ASN:HD21	1.39	0.69
1:B:634:ASN:HD21	1:B:665:ASP:N	1.88	0.69
1:A:634:ASN:HD21	1:A:665:ASP:N	1.91	0.67
1:B:569:LEU:HD12	1:B:570:HIS:H	1.61	0.66
1:A:652:HIS:HB3	1:A:725:ASP:OD1	1.97	0.64
1:A:599:ASN:HB3	1:A:635:GLN:HE22	1.61	0.64
1:A:422:ASP:OD2	1:A:447:HIS:HE1	1.82	0.63
1:A:652:HIS:HA	1:A:724:SER:HB3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:447:HIS:HD2	1:A:455:SER:OG	1.86	0.58
1:B:685:ASN:HD22	1:B:685:ASN:C	2.07	0.57
1:A:517:LYS:HB3	1:A:520:LEU:HD22	1.85	0.57
1:A:605:GLN:O	1:A:606:ASP:OD2	2.22	0.56
1:A:599:ASN:ND2	1:A:632:ASN:H	1.99	0.56
1:A:482:ASN:HD21	1:A:513:LYS:HZ1	1.55	0.55
1:A:414:VAL:HG21	1:A:474:CYS:HA	1.88	0.54
1:A:644:LEU:HD13	1:A:676:PRO:HG2	1.90	0.54
1:A:654:GLY:N	1:A:725:ASP:HB3	2.06	0.54
1:A:669:ARG:NH1	4:A:135:HOH:O	2.40	0.53
1:A:632:ASN:ND2	1:A:634:ASN:H	2.07	0.53
1:A:460:LEU:HD23	1:A:465:TRP:CZ2	2.45	0.52
1:B:632:ASN:HD22	1:B:634:ASN:H	1.55	0.51
1:B:645:HIS:HD2	4:B:104:HOH:O	1.92	0.51
1:B:632:ASN:HD22	1:B:632:ASN:C	2.15	0.50
1:A:632:ASN:C	1:A:632:ASN:HD22	2.15	0.50
1:B:447:HIS:HD2	1:B:455:SER:OG	1.95	0.49
1:B:645:HIS:CD2	4:B:104:HOH:O	2.64	0.49
1:A:511:HIS:HD2	1:A:516:GLU:OE2	1.96	0.49
1:A:686:ALA:O	1:A:690:THR:HG23	2.13	0.48
1:B:416:GLY:HA3	1:B:479:VAL:HG23	1.95	0.47
1:A:482:ASN:HD21	1:A:513:LYS:NZ	2.11	0.47
1:B:422:ASP:OD1	1:B:447:HIS:CE1	2.59	0.47
1:A:476:LEU:HB3	1:A:481:ILE:HD11	1.95	0.47
1:A:525:GLU:HB3	1:A:527:ARG:HH11	1.80	0.46
1:B:590:ASP:O	1:B:594:HIS:HD2	1.99	0.46
1:A:705:GLY:HA2	1:A:706:GLN:HA	1.78	0.46
1:A:489:VAL:HG23	1:A:494:VAL:HG23	1.98	0.46
1:A:632:ASN:HD22	1:A:633:VAL:N	2.14	0.46
1:A:711:THR:CA	1:A:712:TYR:CB	2.67	0.45
1:A:717:ARG:NH1	1:A:718:ASP:OD2	2.50	0.45
1:B:524:PRO:HD2	1:B:568:SER:O	2.16	0.45
1:B:569:LEU:HD13	1:B:570:HIS:H	1.81	0.45
1:A:580:GLY:HA2	1:A:613:GLN:HG3	2.00	0.44
1:B:633:VAL:HG22	4:B:6:HOH:O	2.18	0.44
1:A:710:GLU:O	1:A:711:THR:C	2.56	0.43
1:A:666:LEU:HD23	1:A:666:LEU:HA	1.73	0.43
1:A:693:ARG:HA	1:A:693:ARG:HD2	1.85	0.43
1:A:476:LEU:HB3	1:A:481:ILE:CD1	2.49	0.42
1:A:598:VAL:HG22	4:A:38:HOH:O	2.18	0.42
1:A:654:GLY:HA3	1:A:725:ASP:CB	2.22	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:627:LEU:CD2	1:B:633:VAL:HG13	2.50	0.41
1:B:408:VAL:O	1:B:412:GLN:HB2	2.20	0.41
1:A:592:LEU:HD21	1:A:598:VAL:HG13	2.02	0.41
1:B:569:LEU:HB3	1:B:570:HIS:CD2	2.56	0.41
1:B:585:LEU:HD11	1:B:621:LEU:HB3	2.03	0.41
1:B:683:THR:HG21	4:B:129:HOH:O	2.20	0.41
1:B:662:ARG:NE	4:B:107:HOH:O	2.54	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	275/390 (70%)	261 (95%)	11 (4%)	3 (1%)	14 15
1	B	267/390 (68%)	261 (98%)	6 (2%)	0	100 100
All	All	542/780 (70%)	522 (96%)	17 (3%)	3 (1%)	25 31

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	712	TYR
1	A	606	ASP
1	A	711	THR

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	220/306 (72%)	206 (94%)	14 (6%)	17 23
1	B	204/306 (67%)	183 (90%)	21 (10%)	7 8
All	All	424/612 (69%)	389 (92%)	35 (8%)	11 14

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	407	VAL
1	A	419	GLN
1	A	424	ARG
1	A	456	LYS
1	A	495	LYS
1	A	522	LYS
1	A	598	VAL
1	A	632	ASN
1	A	644	LEU
1	A	660	LEU
1	A	666	LEU
1	A	690	THR
1	A	706	GLN
1	A	712	TYR
1	B	407	VAL
1	B	453	HIS
1	B	468	GLU
1	B	495	LYS
1	B	520	LEU
1	B	525	GLU
1	B	569	LEU
1	B	574	LEU
1	B	602	ASN
1	B	621	LEU
1	B	632	ASN
1	B	633	VAL
1	B	638	SER
1	B	655	LEU
1	B	660	LEU
1	B	662	ARG
1	B	669	ARG
1	B	677	LEU
1	B	685	ASN
1	B	694	LEU

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Mol	Chain	Res	Type
1	B	696	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	410	GLN
1	A	412	GLN
1	A	419	GLN
1	A	442	GLN
1	A	447	HIS
1	A	478	ASN
1	A	482	ASN
1	A	483	GLN
1	A	511	HIS
1	A	599	ASN
1	A	602	ASN
1	A	629	ASN
1	A	632	ASN
1	A	634	ASN
1	A	635	GLN
1	A	706	GLN
1	B	410	GLN
1	B	419	GLN
1	B	447	HIS
1	B	453	HIS
1	B	478	ASN
1	B	570	HIS
1	B	594	HIS
1	B	599	ASN
1	B	602	ASN
1	B	607	ASN
1	B	629	ASN
1	B	632	ASN
1	B	634	ASN
1	B	635	GLN
1	B	645	HIS
1	B	685	ASN
1	B	704	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)

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)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	SO4	B	202	-	4,4,4	0.20	0	6,6,6	0.20	0
3	SO4	A	201	-	4,4,4	0.17	0	6,6,6	0.16	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.

No monomer is involved in short contacts.

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	281/390 (72%)	0.22	20 (7%)	16	21	27, 44, 108, 155	1 (0%)
1	B	273/390 (70%)	0.07	15 (5%)	25	31	26, 45, 110, 144	1 (0%)
All	All	554/780 (71%)	0.15	35 (6%)	20	25	26, 45, 111, 155	2 (0%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	702	ALA	9.7
1	A	528	GLY	9.5
1	B	567	GLY	8.9
1	A	705	GLY	7.4
1	A	703	ALA	6.3
1	A	604	GLY	5.3
1	A	720	SER	5.3
1	A	719	PHE	5.1
1	B	568	SER	5.0
1	A	723	ALA	4.9
1	A	709	ASP	4.9
1	A	704	GLN	4.8
1	B	704	GLN	4.3
1	A	721	LEU	4.3
1	B	703	ALA	4.2
1	A	707	ALA	3.9
1	B	718	ASP	3.8
1	A	725	ASP	3.7
1	B	708	GLY	3.6
1	A	724	SER	3.6
1	A	710	GLU	3.5
1	B	504	GLN	3.2
1	A	718	ASP	3.1
1	A	527	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	522	LYS	2.7
1	B	702	ALA	2.6
1	A	526	ILE	2.5
1	A	711	THR	2.4
1	A	706	GLN	2.3
1	B	569	LEU	2.2
1	B	701	GLU	2.2
1	B	570	HIS	2.1
1	B	605	GLN	2.1
1	B	406	HIS	2.1
1	B	714	ASP	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
3	SO4	A	201	5/5	0.98	0.12	64,66,68,71	0
3	SO4	B	202	5/5	0.98	0.24	75,76,79,84	0
2	ZN	A	999	1/1	1.00	0.13	37,37,37,37	0
2	ZN	B	999	1/1	1.00	0.11	32,32,32,32	0

6.5 Other polymers (i)

There are no such residues in this entry.