Full wwPDB X-ray Structure Validation Report

Jan 31, 2016 – 06:59 PM GMT

PDB ID	:	1DJS
Title	:	LIGAND-BINDING PORTION OF FIBROBLAST GROWTH FACTOR RECEPTOR 2 IN COMPLEX WITH FGF1
Authors	:	Stauber, D.J.; Digabriele, A.D.; Hendrickson, W.A.
Deposited on	:	1999-12-03
Resolution	:	2.40 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

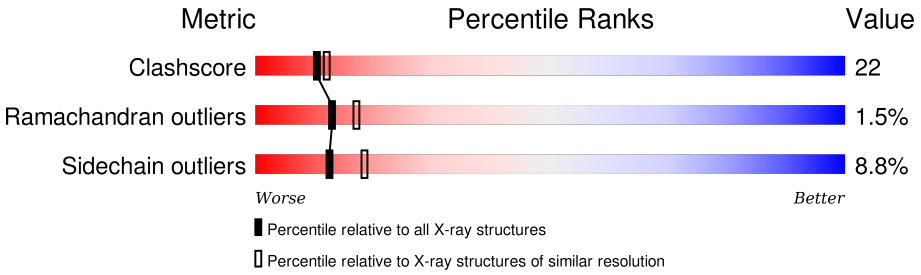
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 2.40 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	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	216	56% 29% 9% • 5%
2	B	135	68% 24% 6% •

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
3	SO4	A	403	-	-	X	-
3	SO4	B	409	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3066 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 PROTEIN (FIBROBLAST GROWTH FACTOR RECEPTOR 2).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	206	1627	1036	284	299	4	4	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	152	GLY	ARG	MUTATION DUE TO SPLICE	UNP P21802
A	162	MSE	MET	MODIFIED	UNP P21802
A	186	MSE	MET	MODIFIED	UNP P21802
A	189	MSE	MET	MODIFIED	UNP P21802
A	218	MSE	MET	MODIFIED	UNP P21802

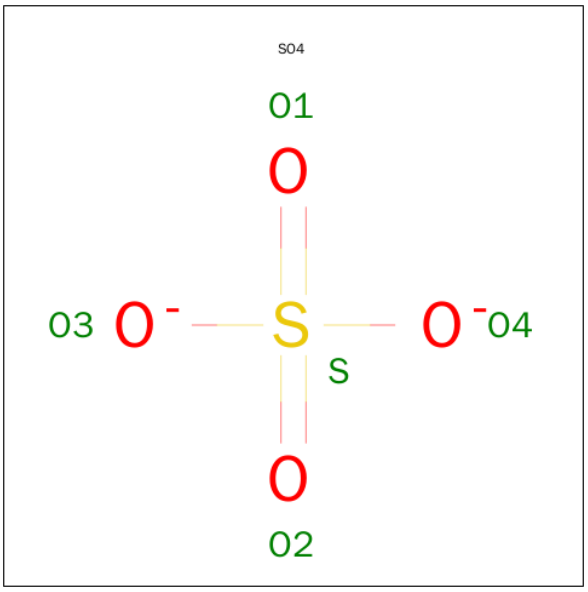
- Molecule 2 is a protein called PROTEIN (FIBROBLAST GROWTH FACTOR 1).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
2	B	135	1075	677	188	206	3	1	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	67	MSE	MET	MODIFIED	UNP P05230

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



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	189	Total O 189 189	0	0
4	B	130	Total O 130 130	0	0

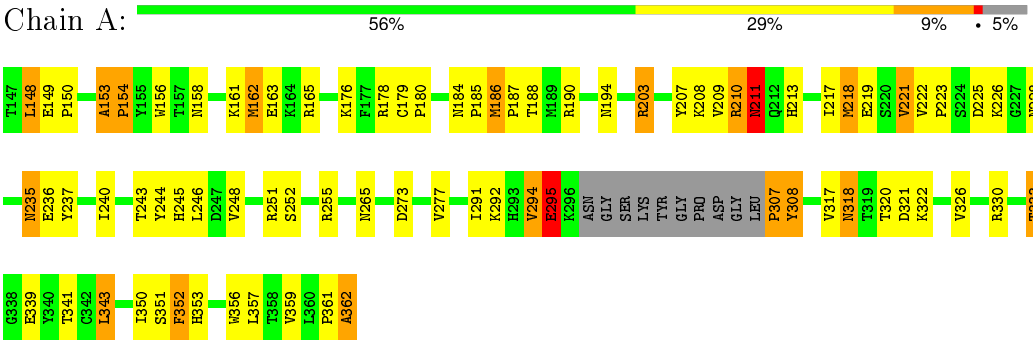
3 Residue-property plots (i)

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.

Note EDS was not executed.

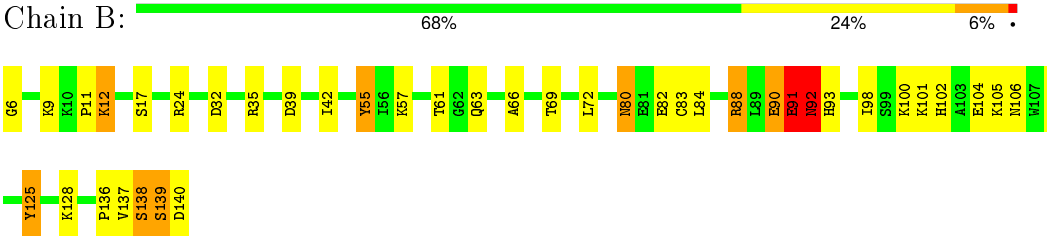
- Molecule 1: PROTEIN (FIBROBLAST GROWTH FACTOR RECEPTOR 2)

Chain A:



- Molecule 2: PROTEIN (FIBROBLAST GROWTH FACTOR 1)

Chain B:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α, β, γ	129.97Å 129.97Å 129.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-2.40)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.221 , 0.315	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3066	wwPDB-VP
Average B, all atoms (Å^2)	30.0	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: 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	0.81	3/1667 (0.2%)	1.38	19/2260 (0.8%)
2	B	0.58	0/1098	1.40	15/1479 (1.0%)
All	All	0.73	3/2765 (0.1%)	1.39	34/3739 (0.9%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	186	MSE	SE-CE	16.39	2.92	1.95
1	A	162	MSE	CG-SE	14.28	2.44	1.95
1	A	218	MSE	CG-SE	8.91	2.25	1.95

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	186	MSE	CG-SE-CE	-15.44	64.94	98.90
1	A	162	MSE	CG-SE-CE	-10.23	76.39	98.90
1	A	218	MSE	CG-SE-CE	-10.23	76.40	98.90
2	B	88	ARG	CD-NE-CZ	10.11	137.76	123.60
2	B	91	GLU	CA-CB-CG	9.91	135.20	113.40
2	B	88	ARG	NE-CZ-NH2	9.49	125.04	120.30
1	A	251	ARG	NE-CZ-NH1	9.37	124.98	120.30
2	B	91	GLU	CB-CG-CD	8.65	137.57	114.20
2	B	91	GLU	C-N-CA	8.64	143.29	121.70
1	A	251	ARG	CD-NE-CZ	8.22	135.11	123.60
1	A	162	MSE	CB-CG-SE	-8.17	88.18	112.70
1	A	218	MSE	CB-CG-SE	-8.03	88.60	112.70
1	A	251	ARG	NE-CZ-NH2	-8.01	116.29	120.30
2	B	91	GLU	O-C-N	-7.84	110.16	122.70
2	B	24	ARG	NE-CZ-NH2	-7.77	116.41	120.30
1	A	153	ALA	CA-C-O	-7.20	104.97	120.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	210	ARG	CD-NE-CZ	7.19	133.66	123.60
2	B	92	ASN	C-N-CA	6.94	139.04	121.70
1	A	154	PRO	N-CA-CB	6.76	111.41	103.30
2	B	90	GLU	O-C-N	-6.53	112.26	122.70
1	A	154	PRO	CA-N-CD	-6.31	102.66	111.50
2	B	55	TYR	CB-CG-CD2	-5.95	117.43	121.00
1	A	308	TYR	N-CA-C	5.88	126.86	111.00
2	B	92	ASN	CA-C-O	5.69	132.05	120.10
2	B	90	GLU	CA-C-N	5.67	129.66	117.20
1	A	330	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	A	362	ALA	N-CA-CB	5.51	117.82	110.10
2	B	91	GLU	CA-C-O	5.45	131.55	120.10
1	A	273	ASP	CB-CG-OD1	-5.26	113.56	118.30
1	A	294	VAL	C-N-CA	5.26	134.85	121.70
1	A	211	ASN	C-N-CA	5.22	134.75	121.70
2	B	92	ASN	N-CA-CB	5.11	119.79	110.60
2	B	32	ASP	CB-CG-OD1	5.09	122.88	118.30
1	A	307	PRO	CA-N-CD	-5.07	104.40	111.50

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	1627	0	1593	74	0
2	B	1075	0	1052	47	0
3	A	20	0	0	3	0
3	B	25	0	0	5	0
4	A	189	0	0	8	0
4	B	130	0	0	12	0
All	All	3066	0	2645	116	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 22.

All (116) 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:218:MSE:CE	1:A:218:MSE:SE	2.19	1.41
1:A:218:MSE:CG	1:A:218:MSE:SE	2.25	1.34
1:A:162:MSE:CG	1:A:162:MSE:SE	2.44	1.16
1:A:218:MSE:SE	1:A:218:MSE:CB	2.68	0.90
2:B:6:GLY:HA3	3:B:409:SO4:O3	1.74	0.87
2:B:136:PRO:HG3	2:B:140:ASP:HA	1.62	0.82
1:A:186:MSE:CE	1:A:186:MSE:HG2	2.09	0.82
2:B:88:ARG:NH1	2:B:98:ILE:HG13	1.95	0.81
1:A:230:THR:OG1	1:A:243:THR:HG22	1.83	0.77
1:A:162:MSE:SE	1:A:162:MSE:CB	2.82	0.77
2:B:92:ASN:N	4:B:410:HOH:O	2.18	0.75
2:B:84:LEU:HD23	2:B:100:LYS:HD2	1.69	0.74
1:A:333:THR:HG22	1:A:335:GLU:H	1.55	0.72
2:B:80:ASN:HD22	2:B:82:GLU:H	1.37	0.70
1:A:222:VAL:HB	1:A:223:PRO:HD2	1.76	0.68
1:A:186:MSE:CE	1:A:186:MSE:SE	2.92	0.68
1:A:218:MSE:SE	1:A:218:MSE:HB3	2.43	0.68
1:A:158:ASN:HD21	1:A:161:LYS:H	1.42	0.67
1:A:190:ARG:HB3	1:A:232:VAL:HG22	1.77	0.67
1:A:318:ASN:HD21	2:B:57:LYS:NZ	1.93	0.65
1:A:186:MSE:CE	1:A:186:MSE:CG	2.74	0.65
1:A:218:MSE:CG	1:A:218:MSE:CE	2.75	0.65
1:A:333:THR:HG22	1:A:335:GLU:N	2.11	0.65
2:B:98:ILE:HD11	2:B:108:PHE:CZ	2.32	0.63
1:A:190:ARG:NH2	1:A:232:VAL:HG21	2.13	0.63
1:A:158:ASN:ND2	1:A:161:LYS:H	1.96	0.63
1:A:162:MSE:CE	1:A:162:MSE:CG	2.76	0.62
2:B:72:LEU:HA	4:B:433:HOH:O	2.01	0.61
2:B:91:GLU:CA	4:B:410:HOH:O	2.47	0.61
2:B:101:LYS:HD3	4:B:428:HOH:O	2.01	0.60
2:B:91:GLU:N	4:B:410:HOH:O	2.36	0.58
2:B:88:ARG:HH11	2:B:98:ILE:HG13	1.69	0.58
1:A:336:ASP:O	1:A:357:LEU:HD23	2.03	0.57
1:A:339:GLU:HG3	3:A:403:SO4:O2	2.03	0.57
1:A:221:VAL:HG13	1:A:248:VAL:HG22	1.86	0.57
2:B:98:ILE:HD11	2:B:108:PHE:CE2	2.39	0.57
1:A:203:ARG:NH2	1:A:219:GLU:O	2.38	0.57
1:A:337:ALA:HB2	1:A:359:VAL:HG23	1.87	0.56
2:B:102:HIS:HA	2:B:104:GLU:OE2	2.05	0.56
1:A:318:ASN:HD21	2:B:57:LYS:HZ2	1.54	0.56
2:B:139:SER:O	2:B:140:ASP:HB2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:136:PRO:CG	2:B:140:ASP:HA	2.33	0.54
1:A:265:ASN:ND2	1:A:356:TRP:H	2.06	0.54
2:B:91:GLU:HB2	4:B:533:HOH:O	2.06	0.54
1:A:317:VAL:HG23	4:B:438:HOH:O	2.07	0.54
1:A:156:TRP:CE2	1:A:240:ILE:HD12	2.42	0.54
1:A:321:ASP:OD1	2:B:9:LYS:HG2	2.07	0.53
2:B:137:VAL:HB	3:B:405:SO4:O3	2.08	0.53
1:A:149:GLU:CB	1:A:150:PRO:HD2	2.38	0.53
1:A:190:ARG:HD2	4:A:555:HOH:O	2.08	0.53
1:A:184:ASN:HA	1:A:185:PRO:C	2.29	0.53
2:B:80:ASN:ND2	2:B:82:GLU:H	2.06	0.53
1:A:162:MSE:SE	1:A:162:MSE:HB2	2.59	0.53
2:B:91:GLU:C	4:B:410:HOH:O	2.47	0.52
1:A:194:ASN:ND2	1:A:228:ASN:H	2.07	0.52
1:A:208:LYS:CD	1:A:210:ARG:HE	2.22	0.52
2:B:11:PRO:HD2	4:B:490:HOH:O	2.10	0.52
1:A:277:VAL:HG22	1:A:326:VAL:HG22	1.92	0.51
2:B:138:SER:O	2:B:140:ASP:N	2.44	0.51
1:A:162:MSE:HG3	1:A:244:TYR:OH	2.11	0.50
1:A:154:PRO:HG2	1:A:240:ILE:HG12	1.94	0.50
1:A:163:GLU:HG3	2:B:35:ARG:NH1	2.26	0.50
2:B:84:LEU:HD23	2:B:100:LYS:CD	2.41	0.50
1:A:333:THR:CG2	1:A:335:GLU:H	2.22	0.50
1:A:176:LYS:HE2	1:A:178:ARG:NH1	2.26	0.50
1:A:356:TRP:HB2	3:A:403:SO4:O4	2.13	0.49
1:A:353:HIS:HB3	4:A:459:HOH:O	2.12	0.49
2:B:80:ASN:HD22	2:B:82:GLU:N	2.09	0.49
1:A:245:HIS:HB3	4:A:505:HOH:O	2.13	0.49
2:B:118:LYS:HE3	3:B:402:SO4:O1	2.13	0.49
2:B:102:HIS:HB3	2:B:105:LYS:HE3	1.94	0.48
1:A:149:GLU:HB3	1:A:150:PRO:HD2	1.95	0.47
1:A:203:ARG:NH1	1:A:225:ASP:OD2	2.47	0.47
2:B:90:GLU:O	2:B:92:ASN:N	2.48	0.47
1:A:221:VAL:HG13	1:A:248:VAL:CG2	2.44	0.47
1:A:235:ASN:ND2	1:A:237:TYR:H	2.13	0.47
1:A:208:LYS:HD3	1:A:210:ARG:HE	1.80	0.46
2:B:72:LEU:HD23	4:B:433:HOH:O	2.15	0.46
2:B:88:ARG:NH2	2:B:125:TYR:OH	2.49	0.45
2:B:69:THR:HG23	4:B:520:HOH:O	2.15	0.45
1:A:158:ASN:HD21	1:A:161:LYS:N	2.12	0.45
1:A:252:SER:HB3	4:A:466:HOH:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:61:THR:OG1	2:B:63:GLN:HG3	2.16	0.45
2:B:39:ASP:O	2:B:42:ILE:HG12	2.17	0.45
2:B:128:LYS:NZ	3:B:408:SO4:O2	2.49	0.44
2:B:136:PRO:HD2	2:B:140:ASP:OD2	2.17	0.44
1:A:333:THR:HG21	1:A:335:GLU:OE1	2.17	0.44
1:A:203:ARG:NH1	1:A:225:ASP:OD1	2.50	0.44
1:A:176:LYS:NZ	3:A:406:SO4:O4	2.51	0.44
1:A:235:ASN:C	1:A:235:ASN:HD22	2.21	0.44
1:A:226:LYS:HE2	4:A:573:HOH:O	2.18	0.44
1:A:320:THR:HB	4:A:563:HOH:O	2.17	0.43
1:A:350:ILE:O	1:A:350:ILE:HG13	2.18	0.43
1:A:211:ASN:HB2	4:A:593:HOH:O	2.18	0.43
1:A:153:ALA:HA	1:A:235:ASN:HD21	1.84	0.43
1:A:207:TYR:CE2	1:A:209:VAL:HG23	2.53	0.43
1:A:186:MSE:HA	1:A:187:PRO:HD3	1.84	0.43
2:B:55:TYR:CZ	2:B:84:LEU:HD13	2.54	0.43
1:A:176:LYS:HD2	1:A:217:ILE:HD11	2.00	0.43
1:A:291:ILE:HG12	1:A:341:THR:HB	2.00	0.43
1:A:317:VAL:HG13	2:B:55:TYR:CD2	2.53	0.43
1:A:361:PRO:O	1:A:362:ALA:CB	2.67	0.42
2:B:91:GLU:HG3	4:B:482:HOH:O	2.18	0.42
2:B:80:ASN:O	2:B:83:CYS:HB2	2.19	0.42
1:A:322:LYS:HG3	4:A:471:HOH:O	2.19	0.42
2:B:91:GLU:O	2:B:93:HIS:CE1	2.72	0.42
2:B:66:ALA:HB2	2:B:83:CYS:SG	2.60	0.41
2:B:105:LYS:O	2:B:106:ASN:HB2	2.20	0.41
1:A:294:VAL:HB	1:A:295:GLU:CD	2.40	0.41
2:B:104:GLU:H	2:B:104:GLU:CD	2.23	0.41
1:A:236:GLU:HG3	1:A:237:TYR:CE2	2.56	0.41
1:A:343:LEU:HD23	1:A:352:PHE:HB3	2.03	0.40
1:A:246:LEU:HD13	1:A:248:VAL:HG23	2.02	0.40
2:B:12:LYS:HD3	3:B:409:SO4:O2	2.22	0.40
1:A:184:ASN:OD1	1:A:185:PRO:HA	2.22	0.40
1:A:179:CYS:N	1:A:180:PRO:CD	2.85	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	202/216 (94%)	190 (94%)	9 (4%)	3 (2%)	13 17
2	B	133/135 (98%)	123 (92%)	8 (6%)	2 (2%)	13 17
All	All	335/351 (95%)	313 (93%)	17 (5%)	5 (2%)	13 17

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	308	TYR
2	B	91	GLU
2	B	139	SER
1	A	295	GLU
1	A	148	LEU

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	176/179 (98%)	158 (90%)	18 (10%)	9 13
2	B	118/117 (101%)	110 (93%)	8 (7%)	20 31
All	All	294/296 (99%)	268 (91%)	26 (9%)	12 18

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	148	LEU

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Mol	Chain	Res	Type
1	A	165	ARG
1	A	188	THR
1	A	203	ARG
1	A	211	ASN
1	A	213	HIS
1	A	221	VAL
1	A	232	VAL
1	A	235	ASN
1	A	255	ARG
1	A	292	LYS
1	A	295	GLU
1	A	307	PRO
1	A	318	ASN
1	A	333	THR
1	A	343	LEU
1	A	351	SER
1	A	352	PHE
2	B	12	LYS
2	B	17	SER
2	B	80	ASN
2	B	91	GLU
2	B	92	ASN
2	B	117	CYS
2	B	125	TYR
2	B	138	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	ASN
1	A	173	ASN
1	A	194	ASN
1	A	235	ASN
1	A	242	HIS
1	A	265	ASN
1	A	318	ASN
2	B	41	HIS
2	B	45	GLN
2	B	80	ASN

5.3.3 RNA

There are no RNA molecules 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

9 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	SO4	A	403	-	4,4,4	0.64	0	6,6,6	0.23	0
3	SO4	A	404	-	4,4,4	0.72	0	6,6,6	0.23	0
3	SO4	A	406	-	4,4,4	0.83	0	6,6,6	0.12	0
3	SO4	A	407	-	4,4,4	0.80	0	6,6,6	0.24	0
3	SO4	B	401	-	4,4,4	0.74	0	6,6,6	0.46	0
3	SO4	B	402	-	4,4,4	0.74	0	6,6,6	0.20	0
3	SO4	B	405	-	4,4,4	0.76	0	6,6,6	0.30	0
3	SO4	B	408	-	4,4,4	0.85	0	6,6,6	0.31	0
3	SO4	B	409	-	4,4,4	0.49	0	6,6,6	1.10	1 (16%)

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	SO4	A	403	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	A	404	-	-	0/0/0/0	0/0/0/0
3	SO4	A	406	-	-	0/0/0/0	0/0/0/0
3	SO4	A	407	-	-	0/0/0/0	0/0/0/0
3	SO4	B	401	-	-	0/0/0/0	0/0/0/0
3	SO4	B	402	-	-	0/0/0/0	0/0/0/0
3	SO4	B	405	-	-	0/0/0/0	0/0/0/0
3	SO4	B	408	-	-	0/0/0/0	0/0/0/0
3	SO4	B	409	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	409	SO4	O2-S-O1	2.47	117.31	109.50

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	SO4	2	0
3	A	406	SO4	1	0
3	B	402	SO4	1	0
3	B	405	SO4	1	0
3	B	408	SO4	1	0
3	B	409	SO4	2	0

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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.