



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 11:09 pm GMT

PDB ID : 1EVT
Title : CRYSTAL STRUCTURE OF FGF1 IN COMPLEX WITH THE EX-
TRACELLULAR LIGAND BINDING DOMAIN OF FGF RECEPTOR 1
(FGFR1)
Authors : Plotnikov, A.N.; Hubbard, S.R.; Schlessinger, J.; Mohammadi, M.
Deposited on : 2000-04-20
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

<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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

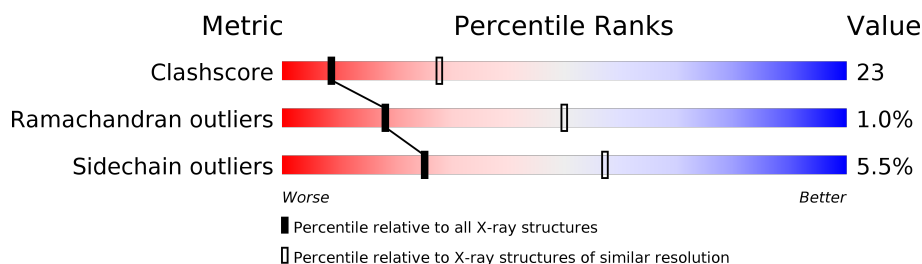
1 Overall quality at a glance

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	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (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 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 $\leq 5\%$. 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	134	
1	B	134	
2	C	225	
2	D	225	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4983 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.

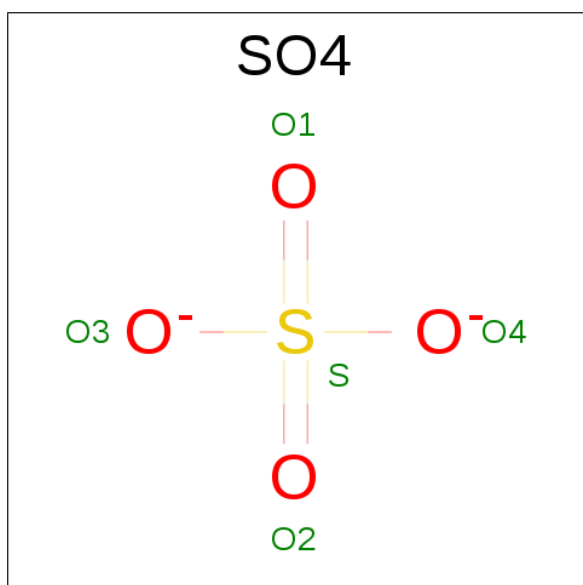
- Molecule 1 is a protein called PROTEIN (FIBROBLAST GROWTH FACTOR 1).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	131	Total	C	N	O	S	0	0	0
			1040	659	181	196	4			
1	B	131	Total	C	N	O	S	0	0	0
			1040	659	181	196	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	191	Total	C	N	O	S	0	0	0
			1439	926	243	262	8			
2	D	192	Total	C	N	O	S	0	0	0
			1444	929	244	263	8			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



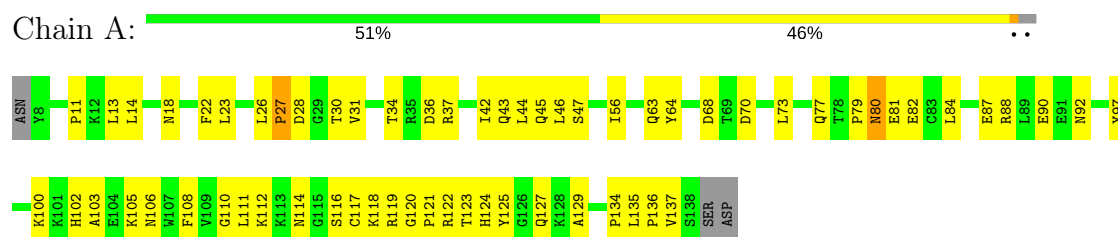
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

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 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.

Note EDS was not executed.

• Molecule 1: PROTEIN (FIBROBLAST GROWTH FACTOR 1)



V325	V322	THR
L326	E233	ASN
H327		ASN
L328	S238	THR
R329		LVS
H330	L245	PRD
V331	D246	N147
S332	V247	
		V151
D335	R254	A152
		P153
Y339	L257	Y154
		W155
L342	K265	T156
A343	T266	S157
		P158
I347	V273	
G348	E274	M161
L349	F275	
S350		V168
H351	V279	P169
H352		A170
	D282	A171
	P283	
		V174
S355	I287	K175
L356	Q288	F176
T357	W289	K177
V358	L290	G178
L359	K291	P179
ALA	H292	S180
LEU	I293	S181
GLU	GLU	T183
ARG	VAL	O184
	ASN	P184
	GLY	N185
	SER	P186
	LVS	T187
	ILE	L188
	GLY	
	PRD	K198
	ASP	P199
	ASN	D200
	LEU	H201
	P306	R202
	Q309	I203
	I310	G204
		G205
		V206
	A314	
	GLY	W213
	VAL	S214
	ASN	
	THR	D218
	THR	S219
	ASP	V220
	LVS	V221
	GLU	
	MET	D224
	E224	

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	62.55Å 64.06Å 64.14Å 93.40° 111.17° 97.18°	Depositor
Resolution (Å)	25.00 – 2.80	Depositor
% Data completeness (in resolution range)	95.7 (25.00-2.80)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.249 , 0.300	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4983	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

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

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	0.46	0/1064	0.69	0/1437
1	B	0.43	0/1064	0.69	1/1437 (0.1%)
2	C	0.47	0/1483	0.69	0/2038
2	D	0.46	0/1488	0.68	0/2043
All	All	0.46	0/5099	0.68	1/6955 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	109	VAL	N-CA-C	-5.15	97.09	111.00

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	1040	0	1014	52	0
1	B	1040	0	1014	43	0
2	C	1439	0	1337	69	0
2	D	1444	0	1346	62	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4983	0	4711	220	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 23.

All (220) 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)
2:D:293:ILE:HD11	2:D:309:GLN:HB2	1.40	1.00
2:D:273:VAL:HG13	2:D:328:LEU:HB2	1.44	0.97
1:B:80:ASN:HD22	1:B:82:GLU:H	1.08	0.94
2:C:293:ILE:HD11	2:C:309:GLN:HB2	1.50	0.93
1:A:27:PRO:HA	1:A:63:GLN:HE22	1.40	0.85
2:C:273:VAL:HG13	2:C:328:LEU:HB2	1.58	0.85
2:C:331:VAL:HG13	2:C:335:ASP:HB2	1.62	0.81
1:A:27:PRO:HA	1:A:63:GLN:NE2	1.96	0.80
2:D:288:GLN:NE2	2:D:310:ILE:HD12	1.98	0.79
1:A:119:ARG:HB2	1:A:122:ARG:HG2	1.62	0.79
1:B:123:THR:HA	1:B:127:GLN:OE1	1.84	0.78
1:A:80:ASN:HD22	1:A:82:GLU:H	1.32	0.77
1:B:119:ARG:HB2	1:B:122:ARG:HG2	1.67	0.77
1:B:80:ASN:HD22	1:B:82:GLU:N	1.83	0.76
1:A:134:PRO:O	1:A:135:LEU:HD23	1.85	0.75
1:B:80:ASN:ND2	1:B:82:GLU:H	1.83	0.75
2:C:332:SER:N	2:C:335:ASP:OD2	2.19	0.75
1:B:27:PRO:HA	1:B:63:GLN:HE22	1.53	0.73
2:C:190:TRP:CH2	2:C:230:CYS:HB3	2.24	0.72
2:C:269:LEU:HD11	2:C:333:PHE:CE1	2.26	0.71
2:D:155:TRP:HB2	2:D:158:PRO:HG3	1.72	0.70
1:B:80:ASN:HD21	1:B:82:GLU:HB2	1.54	0.70
2:C:290:LEU:HD22	2:C:310:ILE:HD13	1.73	0.70
1:B:27:PRO:HA	1:B:63:GLN:NE2	2.09	0.68
1:B:26:LEU:HD12	1:B:30:THR:CG2	2.24	0.67
1:B:68:ASP:OD1	1:B:72:LEU:HB2	1.92	0.67
2:D:174:VAL:HG11	2:D:245:LEU:HD11	1.76	0.67
1:A:123:THR:HA	1:A:127:GLN:OE1	1.95	0.67
2:D:291:LYS:HD3	2:D:339:TYR:CE2	2.30	0.66
2:C:218:ASP:OD2	2:D:265:LYS:HE3	1.95	0.66
2:C:245:LEU:HD23	2:C:246:ASP:N	2.11	0.66
2:D:183:THR:HA	2:D:184:PRO:C	2.17	0.65
2:D:290:LEU:HD22	2:D:310:ILE:HD13	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:291:LYS:HD3	2:D:339:TYR:CZ	2.32	0.65
2:D:170:ALA:O	2:D:171:ALA:HB3	1.96	0.64
1:B:42:ILE:HG13	1:B:43:GLN:NE2	2.13	0.64
1:A:77:GLN:CD	1:A:77:GLN:H	2.02	0.63
2:D:157:SER:N	2:D:158:PRO:HD3	2.13	0.63
1:B:26:LEU:HD12	1:B:30:THR:HG22	1.81	0.62
1:A:119:ARG:HH11	1:A:122:ARG:HD3	1.64	0.61
1:A:22:PHE:CE1	1:A:37:ARG:HG2	2.36	0.61
2:C:157:SER:N	2:C:158:PRO:HD3	2.16	0.61
2:D:273:VAL:CG1	2:D:331:VAL:HG21	2.30	0.61
1:B:88:ARG:HD3	1:B:90:GLU:OE2	2.00	0.61
2:D:342:LEU:HD23	2:D:343:ALA:N	2.17	0.60
2:C:183:THR:HA	2:C:184:PRO:C	2.21	0.60
1:B:112:LYS:HG3	1:B:118:LYS:HB2	1.83	0.60
2:C:331:VAL:HG13	2:C:335:ASP:CB	2.31	0.59
1:A:23:LEU:HD13	1:A:111:LEU:HD11	1.84	0.59
1:A:80:ASN:HD22	1:A:82:GLU:N	1.99	0.59
2:C:245:LEU:C	2:C:245:LEU:HD23	2.23	0.59
2:D:155:TRP:CB	2:D:158:PRO:HG3	2.32	0.59
2:C:170:ALA:O	2:C:171:ALA:HB3	2.03	0.59
2:C:187:THR:O	2:C:232:VAL:HA	2.01	0.59
1:A:42:ILE:HG13	1:A:43:GLN:NE2	2.18	0.58
1:B:16:CYS:HB2	1:B:132:PHE:CE1	2.39	0.58
2:D:187:THR:O	2:D:232:VAL:HA	2.04	0.58
1:B:134:PRO:O	1:B:135:LEU:HD23	2.04	0.57
1:B:77:GLN:CD	1:B:77:GLN:H	2.08	0.57
2:C:161:MET:CE	2:C:178:CYS:HA	2.34	0.57
2:C:291:LYS:HD3	2:C:339:TYR:CZ	2.39	0.56
2:D:174:VAL:HG11	2:D:245:LEU:CD1	2.35	0.56
2:C:276:MET:HE2	2:D:203:ILE:HD12	1.87	0.56
2:D:288:GLN:HE22	2:D:310:ILE:HD12	1.70	0.56
1:A:112:LYS:HG2	1:A:118:LYS:HD2	1.86	0.56
1:A:31:VAL:HG11	1:A:73:LEU:HB3	1.88	0.56
2:C:205:GLY:O	2:C:206:TYR:HB3	2.05	0.56
2:C:331:VAL:HA	2:C:335:ASP:OD2	2.05	0.56
2:C:233:GLU:HB3	2:C:238:SER:HB2	1.88	0.56
1:A:84:LEU:O	1:A:100:LYS:HB2	2.06	0.55
1:A:119:ARG:NH1	1:A:122:ARG:HD3	2.22	0.55
1:A:124:HIS:HB3	1:A:127:GLN:HG3	1.87	0.55
1:B:23:LEU:HD13	1:B:111:LEU:HD11	1.88	0.55
2:C:161:MET:HE2	2:C:178:CYS:HA	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:GLU:HB2	1:B:92:ASN:OD1	2.07	0.55
2:D:282:ASP:HB3	2:D:283:PRO:CD	2.37	0.54
2:C:155:TRP:HB2	2:C:158:PRO:HG3	1.88	0.54
2:D:331:VAL:HG13	2:D:335:ASP:HB2	1.90	0.54
2:C:276:MET:CE	2:D:203:ILE:HD12	2.37	0.54
1:A:80:ASN:ND2	1:A:82:GLU:H	2.03	0.54
2:D:176:PHE:CZ	2:D:245:LEU:HD12	2.43	0.54
2:C:154:TYR:CE1	2:C:181:SER:HB3	2.43	0.54
2:C:289:TRP:CD2	2:C:326:LEU:HB2	2.43	0.53
2:C:288:GLN:NE2	2:C:310:ILE:HD12	2.23	0.53
2:D:156:THR:HG22	2:D:156:THR:O	2.08	0.53
2:D:289:TRP:CD2	2:D:326:LEU:HB2	2.43	0.53
1:B:112:LYS:HG2	1:B:118:LYS:HD2	1.90	0.53
1:A:112:LYS:HG3	1:A:118:LYS:HB2	1.91	0.52
1:A:119:ARG:HB2	1:A:122:ARG:CG	2.37	0.52
2:D:349:LEU:HD12	2:D:349:LEU:C	2.29	0.52
1:B:46:LEU:HG	1:B:56:ILE:HG12	1.91	0.52
2:C:265:LYS:HE3	2:D:218:ASP:OD2	2.10	0.52
1:A:46:LEU:HG	1:A:56:ILE:HG12	1.92	0.52
1:A:102:HIS:ND1	1:A:105:LYS:HE2	2.25	0.51
2:C:257:LEU:CD2	2:C:279:VAL:HG22	2.40	0.51
2:D:170:ALA:O	2:D:171:ALA:CB	2.58	0.51
1:A:14:LEU:HG	1:A:44:LEU:HD12	1.93	0.51
2:D:233:GLU:HB3	2:D:238:SER:HB2	1.92	0.51
1:B:64:TYR:CE1	1:B:79:PRO:HD3	2.46	0.51
2:C:190:TRP:CZ2	2:C:230:CYS:HB3	2.46	0.51
1:A:64:TYR:CE1	1:A:79:PRO:HD3	2.46	0.50
2:C:269:LEU:HD11	2:C:333:PHE:CZ	2.45	0.50
2:D:347:ILE:N	2:D:347:ILE:HD12	2.27	0.50
2:C:156:THR:HG22	2:C:156:THR:O	2.11	0.50
2:C:198:LYS:O	2:C:201:HIS:HB2	2.11	0.50
1:A:80:ASN:HD21	1:A:82:GLU:HB2	1.77	0.50
2:D:151:VAL:HB	2:D:183:THR:HG23	1.93	0.50
2:D:168:VAL:O	2:D:247:VAL:HA	2.11	0.50
2:C:203:ILE:HD13	2:D:325:VAL:HG11	1.94	0.49
1:A:77:GLN:CD	1:A:77:GLN:N	2.65	0.49
2:D:205:GLY:O	2:D:206:TYR:HB3	2.11	0.49
2:D:174:VAL:CG2	2:D:220:VAL:HG11	2.42	0.49
2:D:176:PHE:O	2:D:214:SER:HA	2.12	0.49
2:C:168:VAL:O	2:C:247:VAL:HA	2.11	0.49
1:A:90:GLU:HB2	1:A:92:ASN:OD1	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:347:ILE:HD12	2:C:347:ILE:N	2.28	0.49
2:D:199:PRO:O	2:D:205:GLY:HA2	2.12	0.49
2:D:351:HIS:O	2:D:352:HIS:HD2	1.95	0.49
1:B:120:GLY:N	1:B:121:PRO:HD2	2.28	0.49
2:C:199:PRO:O	2:C:205:GLY:HA2	2.13	0.49
2:D:157:SER:N	2:D:158:PRO:CD	2.75	0.49
1:B:81:GLU:OE2	1:B:101:LYS:HD2	2.12	0.49
1:B:108:PHE:HD1	1:B:123:THR:OG1	1.96	0.49
1:A:88:ARG:HD3	1:A:90:GLU:OE2	2.13	0.48
2:D:161:MET:HE2	2:D:179:PRO:HD2	1.95	0.48
2:D:233:GLU:HB3	2:D:238:SER:CB	2.43	0.48
1:A:13:LEU:HD13	1:A:22:PHE:CE1	2.48	0.48
2:C:190:TRP:CZ3	2:C:230:CYS:HB3	2.48	0.48
1:A:26:LEU:HD12	1:A:30:THR:CG2	2.44	0.48
2:C:265:LYS:HB3	2:C:265:LYS:HE2	1.72	0.48
1:A:27:PRO:CA	1:A:63:GLN:HE22	2.21	0.48
1:B:31:VAL:HG11	1:B:73:LEU:HB3	1.96	0.48
1:B:124:HIS:HB3	1:B:127:GLN:HG3	1.96	0.47
1:B:80:ASN:ND2	1:B:82:GLU:HB2	2.27	0.47
2:C:282:ASP:HB3	2:C:283:PRO:CD	2.44	0.47
2:D:273:VAL:HG12	2:D:331:VAL:HG21	1.96	0.47
1:B:111:LEU:HD12	1:B:132:PHE:CE1	2.49	0.47
2:C:233:GLU:HB3	2:C:238:SER:CB	2.44	0.47
1:A:103:ALA:O	1:A:106:ASN:N	2.42	0.47
2:C:176:PHE:O	2:C:214:SER:HA	2.13	0.47
1:A:31:VAL:HG11	1:A:73:LEU:CB	2.44	0.47
1:B:80:ASN:ND2	1:B:82:GLU:N	2.53	0.47
1:A:42:ILE:HG13	1:A:43:GLN:HE21	1.80	0.47
1:B:73:LEU:HD21	1:B:109:VAL:HG22	1.96	0.47
2:C:257:LEU:HD22	2:C:279:VAL:HG22	1.96	0.46
2:C:254:ARG:HH11	2:C:254:ARG:CG	2.28	0.46
2:D:221:VAL:O	2:D:224:ASP:HB2	2.16	0.46
1:A:108:PHE:HD1	1:A:123:THR:OG1	1.97	0.46
1:A:68:ASP:OD2	1:A:70:ASP:HB2	2.15	0.46
2:C:273:VAL:CG2	2:C:274:GLU:N	2.79	0.46
2:D:153:PRO:HA	2:D:181:SER:O	2.15	0.46
1:A:13:LEU:HD13	1:A:22:PHE:HE1	1.81	0.46
1:A:11:PRO:HG3	1:A:45:GLN:OE1	2.16	0.46
2:C:292:HIS:CD2	2:C:338:GLU:HG2	2.50	0.46
2:D:206:TYR:C	2:D:206:TYR:CD1	2.89	0.46
1:B:100:LYS:O	1:B:100:LYS:HD3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:151:VAL:HB	2:C:183:THR:H	1.82	0.45
2:C:290:LEU:CD2	2:C:310:ILE:HD13	2.43	0.45
2:C:275:PHE:HZ	2:C:356:LEU:HB2	1.82	0.45
2:C:153:PRO:HA	2:C:181:SER:O	2.16	0.45
2:C:154:TYR:N	2:C:154:TYR:CD1	2.84	0.45
2:C:157:SER:N	2:C:158:PRO:CD	2.79	0.45
1:B:37:ARG:NH2	1:B:136:PRO:O	2.50	0.45
2:D:257:LEU:HB2	2:D:352:HIS:CE1	2.51	0.45
2:C:172:LYS:O	2:C:220:VAL:HG22	2.17	0.45
2:C:291:LYS:HD3	2:C:339:TYR:CE2	2.51	0.45
2:C:349:LEU:C	2:C:349:LEU:HD12	2.37	0.45
2:C:154:TYR:CZ	2:C:181:SER:HB3	2.52	0.44
1:A:37:ARG:NH2	1:A:136:PRO:O	2.50	0.44
2:C:170:ALA:O	2:C:171:ALA:CB	2.65	0.44
1:A:105:LYS:O	1:A:106:ASN:HB2	2.17	0.44
1:B:111:LEU:HD12	1:B:132:PHE:HE1	1.82	0.44
2:D:273:VAL:HG11	2:D:331:VAL:HG21	1.99	0.44
1:B:124:HIS:ND1	1:B:125:TYR:N	2.65	0.44
1:B:108:PHE:HB2	1:B:130:ILE:HG21	2.00	0.44
2:D:198:LYS:O	2:D:201:HIS:HB2	2.18	0.44
1:A:87:GLU:HG3	1:A:97:TYR:CE1	2.53	0.43
1:A:124:HIS:ND1	1:A:125:TYR:N	2.65	0.43
2:C:155:TRP:CB	2:C:158:PRO:HG3	2.48	0.43
2:D:245:LEU:HD23	2:D:246:ASP:N	2.32	0.43
1:A:120:GLY:N	1:A:121:PRO:HD2	2.33	0.43
2:D:154:TYR:CZ	2:D:181:SER:HB3	2.54	0.43
1:A:36:ASP:C	1:A:36:ASP:OD2	2.57	0.43
1:B:105:LYS:O	1:B:106:ASN:HB2	2.19	0.43
1:B:14:LEU:HG	1:B:44:LEU:HD12	2.00	0.43
2:C:221:VAL:HB	2:C:222:PRO:HD2	2.01	0.43
1:B:103:ALA:O	1:B:106:ASN:N	2.44	0.42
1:A:110:GLY:O	1:A:111:LEU:HG	2.19	0.42
2:D:347:ILE:CD1	2:D:347:ILE:N	2.82	0.42
2:C:266:THR:HA	2:C:357:THR:O	2.19	0.42
1:A:22:PHE:N	1:A:22:PHE:CD2	2.87	0.42
2:D:176:PHE:CE2	2:D:245:LEU:HD12	2.54	0.42
1:B:108:PHE:CB	1:B:130:ILE:HG21	2.50	0.42
2:C:206:TYR:C	2:C:206:TYR:CD1	2.93	0.42
2:D:161:MET:CE	2:D:178:CYS:HA	2.50	0.42
2:C:161:MET:HE1	2:C:178:CYS:HA	2.01	0.42
1:A:88:ARG:NH1	1:A:90:GLU:OE1	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:275:PHE:HZ	2:D:356:LEU:HB2	1.84	0.42
2:C:276:MET:CE	2:D:203:ILE:HB	2.50	0.42
1:B:119:ARG:NH1	1:B:122:ARG:HD3	2.35	0.41
1:B:68:ASP:OD2	1:B:72:LEU:N	2.33	0.41
2:D:257:LEU:HD22	2:D:279:VAL:HG22	2.01	0.41
2:C:156:THR:CG2	2:C:156:THR:O	2.67	0.41
2:C:272:ASN:HA	2:C:329:ARG:O	2.20	0.41
1:A:37:ARG:HH21	1:A:137:VAL:C	2.24	0.41
2:D:186:PRO:HG2	2:D:213:TRP:CH2	2.54	0.41
2:D:287:ILE:HD12	2:D:287:ILE:N	2.36	0.41
1:A:124:HIS:H	1:A:127:GLN:CD	2.24	0.41
2:C:273:VAL:HG23	2:C:274:GLU:N	2.36	0.41
2:D:266:THR:HG23	2:D:357:THR:HB	2.03	0.41
2:C:197:PHE:HZ	2:C:202:ARG:HD2	1.86	0.41
1:A:106:ASN:HA	1:A:106:ASN:HD22	1.68	0.40
1:A:18:ASN:HB2	1:A:129:ALA:HA	2.03	0.40
2:C:174:VAL:HG11	2:C:245:LEU:CD1	2.52	0.40
2:D:332:SER:N	2:D:335:ASP:OD2	2.47	0.40
2:D:331:VAL:HA	2:D:335:ASP:OD2	2.22	0.40
2:D:342:LEU:C	2:D:342:LEU:HD23	2.42	0.40
1:A:114:ASN:OD1	1:A:116:SER:HB2	2.21	0.40
1:B:13:LEU:HD13	1:B:22:PHE:CE1	2.57	0.40
2:C:331:VAL:CG1	2:C:332:SER:N	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	129/134 (96%)	120 (93%)	8 (6%)	1 (1%)	22	55
1	B	129/134 (96%)	120 (93%)	8 (6%)	1 (1%)	22	55

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	185/225 (82%)	172 (93%)	11 (6%)	2 (1%)	17	47
2	D	186/225 (83%)	170 (91%)	14 (8%)	2 (1%)	17	47
All	All	629/718 (88%)	582 (92%)	41 (6%)	6 (1%)	18	50

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	329	ARG
2	D	329	ARG
2	D	219	SER
2	C	219	SER
1	A	27	PRO
1	B	27	PRO

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	112/118 (95%)	106 (95%)	6 (5%)	26	58
1	B	112/118 (95%)	107 (96%)	5 (4%)	32	66
2	C	150/198 (76%)	140 (93%)	10 (7%)	19	48
2	D	150/198 (76%)	142 (95%)	8 (5%)	26	59
All	All	524/632 (83%)	495 (94%)	29 (6%)	25	57

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ASP
1	A	34	THR
1	A	47	SER
1	A	80	ASN
1	A	81	GLU
1	A	117	CYS

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Mol	Chain	Res	Type
1	B	28	ASP
1	B	34	THR
1	B	47	SER
1	B	80	ASN
1	B	81	GLU
2	C	147	ASN
2	C	154	TYR
2	C	183	THR
2	C	187	THR
2	C	188	LEU
2	C	200	ASP
2	C	254	ARG
2	C	273	VAL
2	C	350	SER
2	C	355	TRP
2	D	147	ASN
2	D	183	THR
2	D	188	LEU
2	D	200	ASP
2	D	254	ARG
2	D	273	VAL
2	D	350	SER
2	D	355	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	63	GLN
1	A	80	ASN
1	A	106	ASN
1	B	41	HIS
1	B	43	GLN
1	B	63	GLN
1	B	80	ASN
1	B	106	ASN
2	C	241	HIS
2	C	284	GLN
2	C	286	HIS
2	C	288	GLN
2	C	352	HIS
2	D	241	HIS

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Mol	Chain	Res	Type
2	D	284	GLN
2	D	286	HIS
2	D	288	GLN
2	D	327	HIS
2	D	352	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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	4002	-	4,4,4	0.33	0	6,6,6	0.20	0
3	SO4	A	4003	-	4,4,4	0.27	0	6,6,6	0.14	0
3	SO4	B	4000	-	4,4,4	0.32	0	6,6,6	0.13	0
3	SO4	B	4001	-	4,4,4	0.32	0	6,6,6	0.20	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	SO4	A	4002	-	-	0/0/0/0	0/0/0/0
3	SO4	A	4003	-	-	0/0/0/0	0/0/0/0
3	SO4	B	4000	-	-	0/0/0/0	0/0/0/0
3	SO4	B	4001	-	-	0/0/0/0	0/0/0/0

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

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.