



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:50 AM GMT

PDB ID : 2ODF
Title : The crystal structure of gene product Atu2144 from Agrobacterium tumefaciens
Authors : Zhang, R.; Xu, X.; Zheng, H.; Savchenko, A.; Edwards, A.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2006-12-22
Resolution : 1.90 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) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

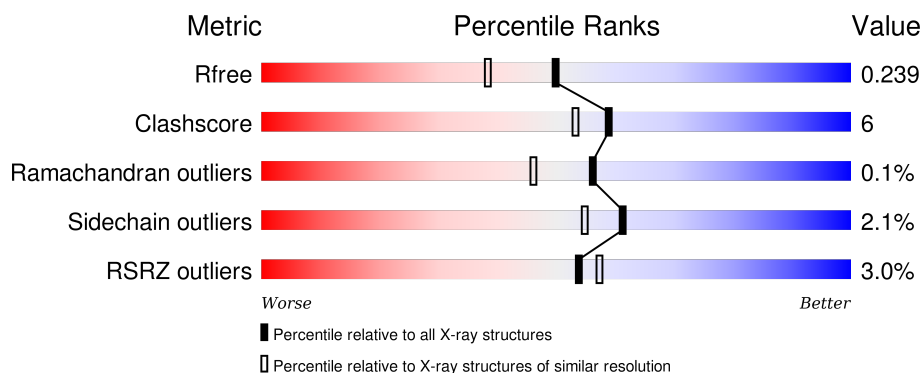
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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.

Mol	Chain	Length	Quality of chain
1	A	257	<div> <div>88%</div> <div>9% ..</div> </div>
1	B	257	<div> <div>2%</div> <div>87%</div> <div>10% ..</div> </div>
1	C	257	<div> <div>%</div> <div>84%</div> <div>11% . .</div> </div>
1	D	257	<div> <div>%</div> <div>87%</div> <div>10% .</div> </div>
1	E	257	<div> <div>3%</div> <div>89%</div> <div>8% .</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	257	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>12%</div> <div>•</div> </div> </div>
1	G	257	<div> <div>7%</div> <div> <div></div> <div>79%</div> <div>16%</div> <div>• •</div> </div> </div>
1	H	257	<div> <div>8%</div> <div> <div></div> <div>83%</div> <div>13%</div> <div>• •</div> </div> </div>

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
2	SO4	A	301	-	X	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17054 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 Hypothetical protein Atu2144.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	0	0	0
			1937	1219	343	369	6			
1	B	252	Total	C	N	O	S	0	0	0
			1928	1215	338	369	6			
1	C	249	Total	C	N	O	S	0	0	0
			1908	1203	337	362	6			
1	D	250	Total	C	N	O	S	0	0	0
			1917	1208	338	365	6			
1	E	251	Total	C	N	O	S	0	0	0
			1928	1214	342	366	6			
1	F	251	Total	C	N	O	S	0	0	0
			1928	1214	342	366	6			
1	G	250	Total	C	N	O	S	0	0	0
			1917	1208	338	365	6			
1	H	251	Total	C	N	O	S	0	0	0
			1928	1214	342	366	6			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	326	Total	O	0	0
			326	326		
3	B	174	Total	O	0	0
			174	174		

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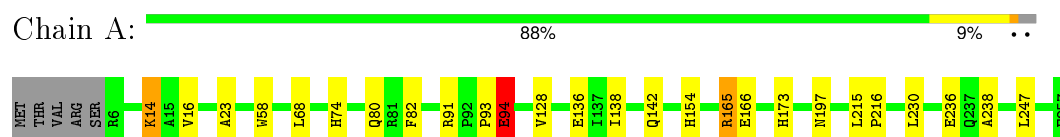
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	216	Total 216	O 216	0	0
3	D	275	Total 275	O 275	0	0
3	E	166	Total 166	O 166	0	0
3	F	171	Total 171	O 171	0	0
3	G	126	Total 126	O 126	0	0
3	H	159	Total 159	O 159	0	0

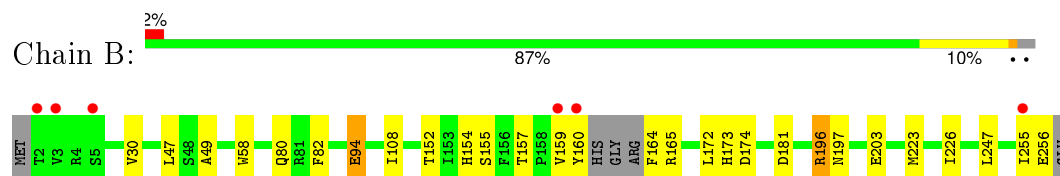
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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.

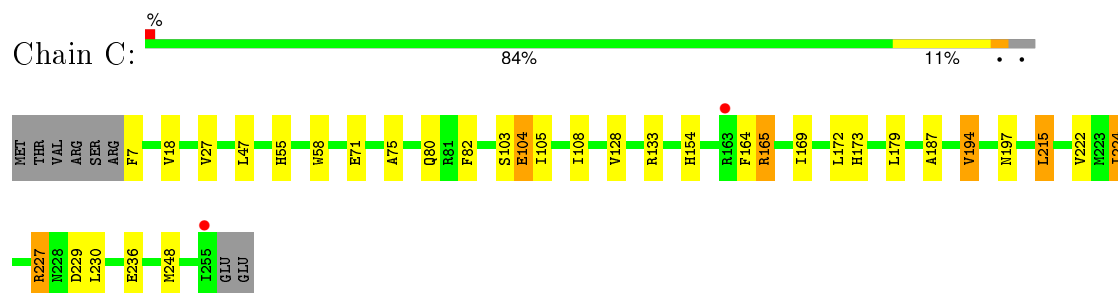
- Molecule 1: Hypothetical protein Atu2144



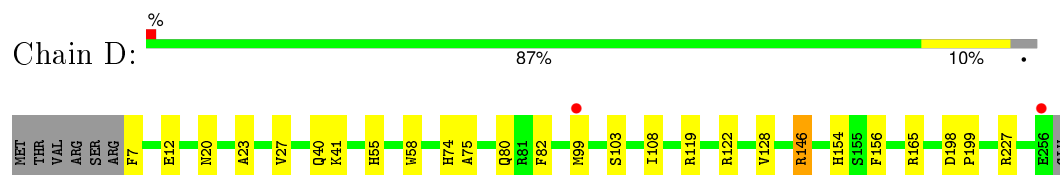
- Molecule 1: Hypothetical protein Atu2144



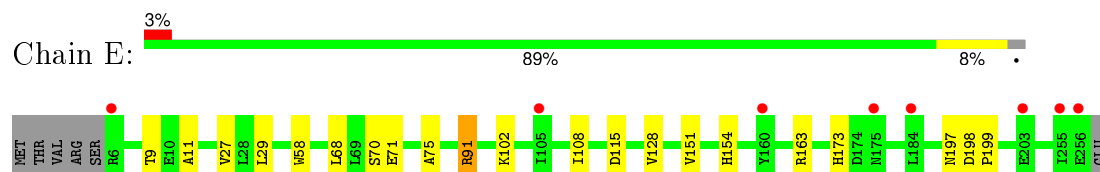
- Molecule 1: Hypothetical protein Atu2144



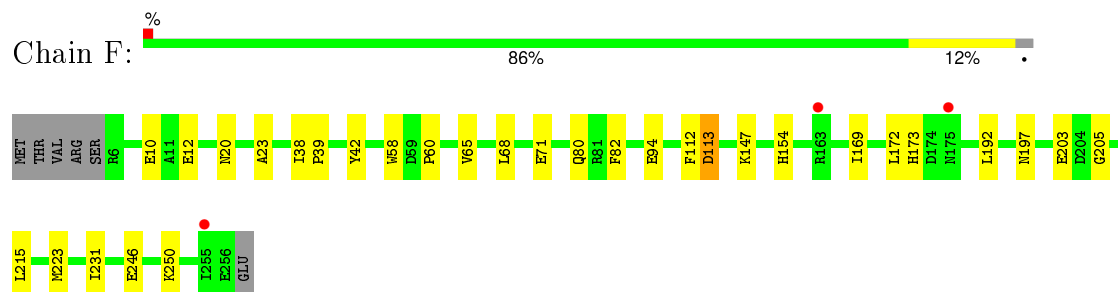
- Molecule 1: Hypothetical protein Atu2144



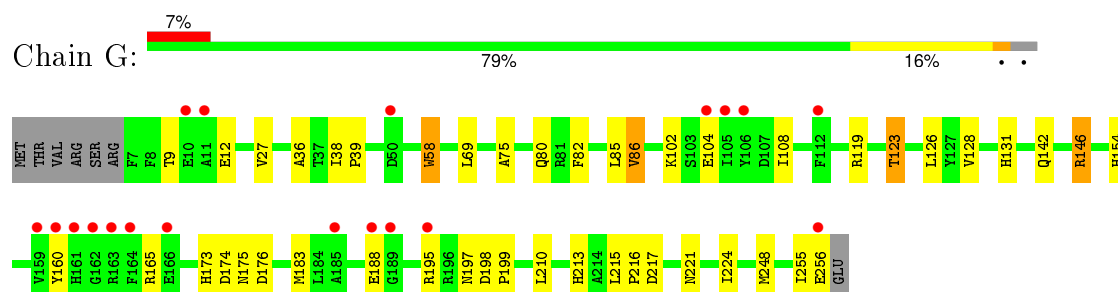
- Molecule 1: Hypothetical protein Atu2144



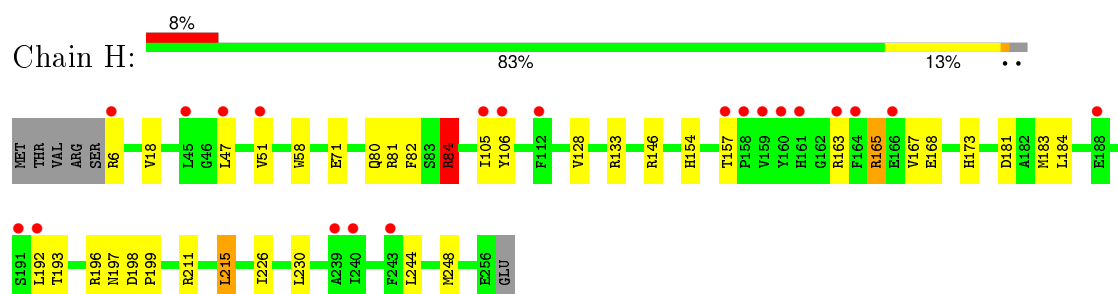
- Molecule 1: Hypothetical protein Atu2144



- Molecule 1: Hypothetical protein Atu2144



- Molecule 1: Hypothetical protein Atu2144



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	110.52Å 89.77Å 120.36Å 90.00° 91.35° 90.00°	Depositor
Resolution (Å)	120.00 – 1.90 45.21 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.0 (120.00-1.90) 98.9 (45.21-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.189 , 0.234 0.194 , 0.239	Depositor DCC
R_{free} test set	9148 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	25.8	Xtriage
Anisotropy	0.500	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 57.8	EDS
Estimated twinning fraction	0.026 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 182396 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17054	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

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	0.76	1/1977 (0.1%)	0.75	3/2684 (0.1%)
1	B	0.61	0/1966	0.69	2/2670 (0.1%)
1	C	0.66	0/1948	0.72	4/2646 (0.2%)
1	D	0.67	0/1957	0.72	2/2658 (0.1%)
1	E	0.55	0/1968	0.67	2/2672 (0.1%)
1	F	0.58	0/1968	0.66	0/2672
1	G	0.55	0/1957	0.65	2/2658 (0.1%)
1	H	0.57	0/1968	0.69	2/2672 (0.1%)
All	All	0.62	1/15709 (0.0%)	0.70	17/21332 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	94	GLU	CB-CG	5.88	1.63	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	165	ARG	NE-CZ-NH1	7.92	124.26	120.30
1	E	91	ARG	NE-CZ-NH2	-6.60	117.00	120.30
1	C	227	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	C	165	ARG	NE-CZ-NH2	-6.42	117.09	120.30
1	D	165	ARG	NE-CZ-NH2	-6.31	117.15	120.30
1	B	196	ARG	NE-CZ-NH2	-6.16	117.22	120.30
1	A	165	ARG	NE-CZ-NH2	-5.89	117.36	120.30
1	G	146	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	D	146	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	A	91	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	A	91	ARG	NE-CZ-NH2	-5.35	117.62	120.30
1	B	196	ARG	NE-CZ-NH1	5.32	122.96	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	84	ARG	NE-CZ-NH1	5.30	122.95	120.30
1	E	115	ASP	CB-CG-OD1	5.22	123.00	118.30
1	G	146	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	C	227	ARG	NE-CZ-NH1	5.02	122.81	120.30
1	H	165	ARG	NE-CZ-NH2	-5.01	117.80	120.30

There are no chirality outliers.

There are no planarity outliers.

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	1937	0	1902	17	0
1	B	1928	0	1895	21	0
1	C	1908	0	1877	38	0
1	D	1917	0	1883	17	0
1	E	1928	0	1896	14	0
1	F	1928	0	1896	20	0
1	G	1917	0	1883	31	0
1	H	1928	0	1896	23	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	10	0	0	1	0
2	D	10	0	0	0	0
2	F	10	0	0	0	0
2	G	5	0	0	0	0
2	H	5	0	0	1	0
3	A	326	0	0	6	0
3	B	174	0	0	5	0
3	C	216	0	0	6	0
3	D	275	0	0	7	0
3	E	166	0	0	5	0
3	F	171	0	0	7	0
3	G	126	0	0	6	0
3	H	159	0	0	3	0
All	All	17054	0	15128	176	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 6.

All (176) 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:136:GLU:OE1	3:A:539:HOH:O	1.75	1.02
1:F:205:GLY:O	3:F:479:HOH:O	1.76	1.01
1:G:224:ILE:HD11	1:G:248:MET:CE	1.96	0.96
1:C:227:ARG:HD2	1:C:229:ASP:OD1	1.69	0.92
1:B:154:HIS:HD2	3:B:303:HOH:O	1.51	0.91
1:C:222:VAL:CG1	1:C:248:MET:HE1	2.01	0.91
1:C:215:LEU:HD13	3:C:512:HOH:O	1.75	0.85
1:C:179:LEU:HD21	1:C:248:MET:HE2	1.58	0.83
1:C:179:LEU:HD21	1:C:248:MET:CE	2.10	0.82
1:F:94:GLU:OE1	3:F:478:HOH:O	1.98	0.82
1:E:29:LEU:HD23	1:E:151:VAL:HG22	1.62	0.81
1:G:119:ARG:O	1:G:123:THR:HG23	1.82	0.80
1:E:102:LYS:HA	1:E:108:ILE:HD12	1.65	0.79
1:E:91:ARG:NE	3:E:423:HOH:O	2.17	0.77
1:G:154:HIS:HD2	3:G:335:HOH:O	1.68	0.76
1:E:68:LEU:O	1:E:71:GLU:HG3	1.87	0.75
1:D:146:ARG:HD2	3:D:360:HOH:O	1.85	0.75
1:B:172:LEU:HD22	1:B:197:ASN:HA	1.68	0.74
1:D:154:HIS:HD2	3:D:309:HOH:O	1.71	0.73
1:G:224:ILE:HD11	1:G:248:MET:HE1	1.69	0.72
1:H:168:GLU:HG3	1:H:193:THR:HB	1.72	0.72
1:C:173:HIS:H	1:C:197:ASN:HD21	1.38	0.71
1:B:173:HIS:H	1:B:197:ASN:HD21	1.38	0.70
1:E:9:THR:HG21	3:E:384:HOH:O	1.90	0.70
1:C:80:GLN:HE21	1:C:82:PHE:H	1.39	0.70
1:H:154:HIS:HD2	3:H:367:HOH:O	1.74	0.70
1:G:224:ILE:HD11	1:G:248:MET:HE3	1.74	0.69
1:D:99:MET:HE3	1:D:119:ARG:HG2	1.73	0.69
1:B:47:LEU:HD21	1:B:108:ILE:HD13	1.73	0.69
1:C:222:VAL:HG12	1:C:248:MET:HE1	1.73	0.69
1:F:223:MET:SD	3:F:443:HOH:O	2.51	0.68
1:G:80:GLN:HE21	1:G:82:PHE:H	1.40	0.67
1:A:173:HIS:H	1:A:197:ASN:HD21	1.41	0.67
1:A:80:GLN:HE21	1:A:82:PHE:H	1.41	0.67
1:H:71:GLU:HG2	3:H:438:HOH:O	1.96	0.66
1:D:80:GLN:HE21	1:D:82:PHE:H	1.40	0.66
1:C:222:VAL:HG11	1:C:248:MET:HE1	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:131:HIS:ND1	1:G:213:HIS:HE1	1.93	0.66
1:G:146:ARG:HD2	3:G:317:HOH:O	1.97	0.65
1:F:215:LEU:HD22	3:F:472:HOH:O	1.98	0.64
1:C:222:VAL:CG1	1:C:248:MET:CE	2.74	0.64
1:E:128:VAL:HG21	1:G:128:VAL:HG21	1.79	0.63
1:F:80:GLN:HE21	1:F:82:PHE:H	1.44	0.62
1:D:41:LYS:HE3	3:D:524:HOH:O	2.00	0.62
1:E:173:HIS:H	1:E:197:ASN:HD21	1.47	0.62
1:C:227:ARG:CD	1:C:229:ASP:OD1	2.44	0.61
1:A:154:HIS:HD2	3:A:309:HOH:O	1.81	0.61
1:C:222:VAL:HG11	1:C:248:MET:CE	2.31	0.60
1:H:168:GLU:HB3	1:H:230:LEU:HD11	1.82	0.60
1:C:222:VAL:HG13	1:C:224:ILE:CD1	2.32	0.59
1:D:41:LYS:NZ	3:D:487:HOH:O	2.35	0.59
1:G:173:HIS:H	1:G:197:ASN:HD21	1.48	0.59
3:E:300:HOH:O	1:G:123:THR:HG21	2.03	0.59
1:B:80:GLN:HE21	1:B:82:PHE:H	1.49	0.59
1:F:68:LEU:O	1:F:71:GLU:HG3	2.04	0.58
1:H:173:HIS:H	1:H:197:ASN:HD21	1.51	0.58
1:A:166:GLU:HG2	3:A:590:HOH:O	2.04	0.57
1:C:47:LEU:HD21	1:C:108:ILE:HD13	1.85	0.57
1:F:173:HIS:H	1:F:197:ASN:HD21	1.54	0.56
1:A:14:LYS:HD2	3:A:347:HOH:O	2.06	0.56
1:F:60:PRO:HD2	3:F:313:HOH:O	2.06	0.55
1:B:173:HIS:HE1	1:B:196:ARG:HE	1.53	0.55
1:B:181:ASP:OD1	1:B:196:ARG:NH2	2.36	0.55
1:F:112:PHE:O	1:F:113:ASP:C	2.44	0.55
1:E:154:HIS:HD2	3:E:264:HOH:O	1.90	0.55
1:H:226:ILE:HD11	1:H:244:LEU:HD11	1.89	0.55
1:G:213:HIS:HD2	3:G:311:HOH:O	1.90	0.55
1:H:183:MET:SD	1:H:248:MET:HE2	2.47	0.54
1:H:84:ARG:HD3	3:H:339:HOH:O	2.07	0.54
1:H:157:THR:HG22	1:H:165:ARG:HH21	1.73	0.54
1:B:173:HIS:HD2	1:B:174:ASP:O	1.90	0.54
1:H:184:LEU:CD1	1:H:196:ARG:HB2	2.38	0.53
1:C:7:PHE:N	3:C:410:HOH:O	2.41	0.53
1:B:247:LEU:HD12	3:B:325:HOH:O	2.07	0.53
1:F:154:HIS:HD2	3:F:328:HOH:O	1.91	0.53
1:E:9:THR:HG23	3:E:286:HOH:O	2.08	0.53
1:C:154:HIS:HE1	3:C:306:HOH:O	1.92	0.53
1:C:222:VAL:HG13	1:C:224:ILE:HD11	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:LYS:HG3	1:A:16:VAL:H	1.74	0.53
1:A:23:ALA:HB1	1:A:74:HIS:O	2.08	0.53
1:D:41:LYS:CE	3:D:524:HOH:O	2.54	0.53
1:F:147:LYS:NZ	3:F:424:HOH:O	2.40	0.53
1:C:187:ALA:HB1	1:C:194:VAL:HG13	1.91	0.53
1:C:165:ARG:HD3	3:C:423:HOH:O	2.08	0.52
1:C:179:LEU:HD21	1:C:248:MET:HE3	1.89	0.52
1:G:210:LEU:HD13	1:G:221:ASN:HD22	1.75	0.51
1:H:157:THR:HG22	1:H:165:ARG:NH2	2.25	0.51
1:H:47:LEU:HD22	1:H:51:VAL:HG11	1.93	0.51
1:D:7:PHE:N	3:D:547:HOH:O	2.44	0.51
1:H:181:ASP:OD1	1:H:196:ARG:NH2	2.43	0.51
1:H:81:ARG:NH1	2:H:310:SO4:O2	2.44	0.50
1:H:184:LEU:HD13	1:H:196:ARG:HB2	1.93	0.50
1:B:94:GLU:OE2	3:B:322:HOH:O	2.19	0.50
1:H:18:VAL:HG21	1:H:133:ARG:CZ	2.41	0.50
1:D:198:ASP:OD1	1:D:199:PRO:HA	2.12	0.49
1:D:99:MET:SD	1:D:122:ARG:HD2	2.52	0.49
1:A:165:ARG:HD3	3:A:504:HOH:O	2.11	0.49
1:C:222:VAL:HG12	1:C:248:MET:CE	2.38	0.49
1:G:142:GLN:HE22	1:G:217:ASP:CG	2.15	0.49
1:F:65:VAL:HG21	1:F:231:ILE:HG21	1.95	0.49
1:D:12:GLU:OE2	1:D:40:GLN:N	2.44	0.49
1:E:27:VAL:HG13	1:E:151:VAL:HG13	1.94	0.48
1:G:86:VAL:HG22	1:G:126:LEU:CD1	2.44	0.48
1:B:159:VAL:O	1:B:160:TYR:HB2	2.14	0.48
1:F:203:GLU:H	1:F:203:GLU:CD	2.17	0.48
1:C:104:GLU:OE2	2:C:303:SO4:O3	2.30	0.48
1:C:128:VAL:HG21	1:H:128:VAL:HG21	1.95	0.48
1:D:103:SER:HB2	1:D:108:ILE:HD11	1.96	0.48
1:B:49:ALA:HB3	1:C:164:PHE:HB3	1.96	0.48
1:G:176:ASP:N	3:G:398:HOH:O	2.46	0.47
1:G:38:ILE:HG22	1:G:39:PRO:O	2.13	0.47
1:F:246:GLU:HG3	1:F:250:LYS:NZ	2.29	0.47
1:A:93:PRO:HD2	1:A:94:GLU:OE1	2.14	0.47
1:C:105:ILE:O	1:C:105:ILE:HG22	2.13	0.47
1:H:80:GLN:HE21	1:H:82:PHE:H	1.62	0.47
1:B:255:ILE:O	1:B:256:GLU:CB	2.63	0.47
1:E:9:THR:HG22	1:E:11:ALA:N	2.28	0.47
1:A:247:LEU:HD13	3:A:496:HOH:O	2.15	0.47
1:G:27:VAL:O	1:G:75:ALA:HB1	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:160:TYR:C	3:B:352:HOH:O	2.54	0.46
1:E:68:LEU:O	1:E:71:GLU:CG	2.62	0.46
1:A:68:LEU:HD13	1:A:238:ALA:HA	1.97	0.46
1:C:103:SER:O	1:C:105:ILE:N	2.49	0.46
1:C:230:LEU:O	1:C:236:GLU:HB3	2.16	0.46
1:G:195:ARG:NE	3:G:403:HOH:O	2.48	0.45
1:C:103:SER:O	1:C:103:SER:OG	2.30	0.45
1:D:156:PHE:CE2	1:D:227:ARG:HG3	2.50	0.45
1:G:255:ILE:O	1:G:256:GLU:C	2.54	0.45
1:C:222:VAL:CG1	1:C:224:ILE:HD11	2.47	0.45
1:A:230:LEU:O	1:A:236:GLU:HB3	2.17	0.45
1:H:81:ARG:HH11	1:H:81:ARG:HG2	1.81	0.45
1:A:14:LYS:HE2	1:A:16:VAL:H	1.82	0.44
1:A:138:ILE:O	1:A:142:GLN:HG3	2.17	0.44
1:B:164:PHE:O	1:B:165:ARG:CB	2.65	0.44
1:H:168:GLU:OE1	1:H:192:LEU:HA	2.17	0.44
1:C:169:ILE:O	1:C:194:VAL:HA	2.17	0.44
1:F:20:ASN:HB3	1:F:23:ALA:HB2	1.99	0.44
1:G:80:GLN:HE21	1:G:82:PHE:N	2.13	0.44
1:F:172:LEU:HG	1:F:197:ASN:HD22	1.82	0.44
1:H:211:ARG:HA	1:H:215:LEU:HD22	2.00	0.44
1:A:128:VAL:HG21	1:D:128:VAL:HG21	1.99	0.44
1:B:154:HIS:CD2	3:B:303:HOH:O	2.42	0.44
1:B:203:GLU:H	1:B:203:GLU:CD	2.21	0.44
1:G:160:TYR:HB2	1:G:165:ARG:CZ	2.48	0.44
1:G:215:LEU:HD21	3:G:433:HOH:O	2.18	0.44
1:G:215:LEU:N	1:G:216:PRO:CD	2.81	0.43
1:F:80:GLN:HE21	1:F:82:PHE:N	2.14	0.43
1:E:198:ASP:OD1	1:E:199:PRO:HA	2.18	0.43
1:G:210:LEU:HD13	1:G:221:ASN:ND2	2.34	0.43
1:H:105:ILE:HD11	1:H:106:TYR:CZ	2.53	0.43
1:F:169:ILE:HD11	1:F:192:LEU:HD13	2.01	0.43
1:C:18:VAL:HG21	1:C:133:ARG:CZ	2.49	0.43
1:G:36:ALA:HB2	1:G:58:TRP:CG	2.54	0.42
1:G:9:THR:N	1:G:12:GLU:OE1	2.43	0.42
1:G:198:ASP:OD1	1:G:199:PRO:HA	2.19	0.42
1:G:119:ARG:O	1:G:123:THR:CG2	2.62	0.42
1:C:7:PHE:N	3:C:493:HOH:O	2.51	0.42
1:B:172:LEU:HB2	1:B:223:MET:HB3	2.01	0.42
1:G:85:LEU:HD22	1:G:108:ILE:CD1	2.50	0.42
1:D:55:HIS:HD2	3:D:396:HOH:O	2.03	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:55:HIS:HE1	1:C:104:GLU:OE1	2.02	0.41
1:C:27:VAL:O	1:C:75:ALA:HB1	2.20	0.41
1:G:183:MET:SD	1:G:248:MET:HE2	2.60	0.41
1:C:224:ILE:N	1:C:224:ILE:CD1	2.84	0.41
1:C:215:LEU:CD1	3:C:512:HOH:O	2.49	0.41
1:B:49:ALA:HB3	1:C:164:PHE:CB	2.50	0.41
1:B:30:VAL:HB	1:B:152:THR:HG22	2.02	0.41
1:D:27:VAL:O	1:D:75:ALA:HB1	2.20	0.41
1:F:12:GLU:OE2	1:F:39:PRO:HA	2.21	0.41
1:C:172:LEU:HA	1:C:197:ASN:ND2	2.36	0.41
1:D:23:ALA:HB1	1:D:74:HIS:O	2.21	0.41
1:H:198:ASP:OD1	1:H:199:PRO:HA	2.21	0.41
1:B:155:SER:HA	1:B:226:ILE:O	2.20	0.40
1:A:215:LEU:N	1:A:216:PRO:CD	2.84	0.40
1:F:38:ILE:HG22	1:F:42:TYR:HB2	2.02	0.40
1:E:70:SER:OG	1:E:75:ALA:O	2.28	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	250/257 (97%)	245 (98%)	5 (2%)	0	100	100
1	B	248/257 (96%)	244 (98%)	4 (2%)	0	100	100
1	C	247/257 (96%)	239 (97%)	7 (3%)	1 (0%)	39	27
1	D	248/257 (96%)	243 (98%)	5 (2%)	0	100	100
1	E	249/257 (97%)	243 (98%)	6 (2%)	0	100	100
1	F	249/257 (97%)	238 (96%)	10 (4%)	1 (0%)	39	27
1	G	248/257 (96%)	241 (97%)	7 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	249/257 (97%)	244 (98%)	5 (2%)	0	100	100
All	All	1988/2056 (97%)	1937 (97%)	49 (2%)	2 (0%)	56	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	104	GLU
1	F	113	ASP

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	202/207 (98%)	199 (98%)	3 (2%)	72	69
1	B	202/207 (98%)	199 (98%)	3 (2%)	72	69
1	C	199/207 (96%)	194 (98%)	5 (2%)	55	47
1	D	200/207 (97%)	198 (99%)	2 (1%)	82	81
1	E	201/207 (97%)	199 (99%)	2 (1%)	82	81
1	F	201/207 (97%)	199 (99%)	2 (1%)	82	81
1	G	200/207 (97%)	191 (96%)	9 (4%)	34	21
1	H	201/207 (97%)	194 (96%)	7 (4%)	43	31
All	All	1606/1656 (97%)	1573 (98%)	33 (2%)	61	55

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LYS
1	A	58	TRP
1	A	94	GLU
1	B	58	TRP
1	B	94	GLU
1	B	157	THR

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Mol	Chain	Res	Type
1	C	58	TRP
1	C	71	GLU
1	C	194	VAL
1	C	215	LEU
1	C	224	ILE
1	D	20	ASN
1	D	58	TRP
1	E	58	TRP
1	E	163	ARG
1	F	10	GLU
1	F	58	TRP
1	G	58	TRP
1	G	69	LEU
1	G	86	VAL
1	G	102	LYS
1	G	104	GLU
1	G	123	THR
1	G	174	ASP
1	G	175	ASN
1	G	188	GLU
1	H	6	ARG
1	H	58	TRP
1	H	84	ARG
1	H	146	ARG
1	H	163	ARG
1	H	167	VAL
1	H	215	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	HIS
1	A	80	GLN
1	A	154	HIS
1	A	197	ASN
1	B	80	GLN
1	B	154	HIS
1	B	173	HIS
1	B	197	ASN
1	B	221	ASN
1	C	80	GLN
1	C	154	HIS

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Mol	Chain	Res	Type
1	C	175	ASN
1	C	197	ASN
1	C	221	ASN
1	D	55	HIS
1	D	80	GLN
1	D	154	HIS
1	D	161	HIS
1	D	221	ASN
1	E	20	ASN
1	E	154	HIS
1	E	161	HIS
1	E	197	ASN
1	E	245	HIS
1	F	20	ASN
1	F	80	GLN
1	F	154	HIS
1	F	161	HIS
1	F	175	ASN
1	F	197	ASN
1	G	80	GLN
1	G	154	HIS
1	G	197	ASN
1	G	213	HIS
1	G	221	ASN
1	H	20	ASN
1	H	80	GLN
1	H	154	HIS
1	H	197	ASN
1	H	221	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 ⓘ

10 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	301	-	4,4,4	3.60	4 (100%)	6,6,6	0.80	0
2	SO4	B	302	-	4,4,4	0.28	0	6,6,6	0.42	0
2	SO4	C	303	-	4,4,4	0.29	0	6,6,6	0.40	0
2	SO4	C	304	-	4,4,4	0.39	0	6,6,6	0.16	0
2	SO4	D	305	-	4,4,4	0.51	0	6,6,6	0.41	0
2	SO4	D	306	-	4,4,4	0.55	0	6,6,6	0.48	0
2	SO4	F	307	-	4,4,4	0.41	0	6,6,6	0.27	0
2	SO4	F	308	-	4,4,4	0.46	0	6,6,6	0.75	0
2	SO4	G	309	-	4,4,4	0.21	0	6,6,6	0.12	0
2	SO4	H	310	-	4,4,4	0.35	0	6,6,6	0.29	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
2	SO4	A	301	-	-	0/0/0/0	0/0/0/0
2	SO4	B	302	-	-	0/0/0/0	0/0/0/0
2	SO4	C	303	-	-	0/0/0/0	0/0/0/0
2	SO4	C	304	-	-	0/0/0/0	0/0/0/0
2	SO4	D	305	-	-	0/0/0/0	0/0/0/0
2	SO4	D	306	-	-	0/0/0/0	0/0/0/0
2	SO4	F	307	-	-	0/0/0/0	0/0/0/0
2	SO4	F	308	-	-	0/0/0/0	0/0/0/0
2	SO4	G	309	-	-	0/0/0/0	0/0/0/0
2	SO4	H	310	-	-	0/0/0/0	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	SO4	O2-S	-4.29	1.32	1.47
2	A	301	SO4	O4-S	-3.68	1.34	1.47
2	A	301	SO4	O3-S	-3.37	1.35	1.47
2	A	301	SO4	O1-S	-2.93	1.37	1.47

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	303	SO4	1	0
2	H	310	SO4	1	0

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 ⓘ

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	252/257 (98%)	-0.13	0	100	100	15, 20, 32, 44	0
1	B	252/257 (98%)	0.10	6 (2%)	62	66	18, 28, 43, 57	0
1	C	249/257 (96%)	0.02	2 (0%)	87	88	18, 27, 36, 46	0
1	D	250/257 (97%)	-0.04	2 (0%)	87	88	15, 24, 33, 48	0
1	E	251/257 (97%)	0.16	8 (3%)	51	54	20, 31, 44, 50	0
1	F	251/257 (97%)	0.03	3 (1%)	81	83	18, 29, 40, 51	0
1	G	250/257 (97%)	0.50	19 (7%)	17	18	21, 33, 53, 63	0
1	H	251/257 (97%)	0.53	21 (8%)	14	15	17, 33, 56, 68	0
All	All	2006/2056 (97%)	0.15	61 (3%)	54	57	15, 28, 47, 68	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	105	ILE	8.2
1	G	106	TYR	7.4
1	G	160	TYR	5.7
1	H	164	PHE	5.4
1	E	105	ILE	5.3
1	H	159	VAL	4.9
1	H	105	ILE	4.8
1	H	106	TYR	4.7
1	C	163	ARG	4.7
1	G	159	VAL	4.4
1	G	11	ALA	4.4
1	H	112	PHE	4.2
1	G	164	PHE	4.1
1	H	166	GLU	3.9
1	H	163	ARG	3.8
1	G	163	ARG	3.7

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Mol	Chain	Res	Type	RSRZ
1	G	189	GLY	3.6
1	G	188	GLU	3.5
1	F	163	ARG	3.5
1	H	160	TYR	3.5
1	E	255	ILE	3.4
1	H	239	ALA	3.4
1	G	166	GLU	3.4
1	H	6	ARG	3.4
1	B	5	SER	3.3
1	H	243	PHE	3.3
1	H	158	PRO	3.2
1	H	161	HIS	3.2
1	H	188	GLU	3.0
1	B	2	THR	3.0
1	B	3	VAL	2.8
1	G	104	GLU	2.7
1	G	185	ALA	2.7
1	E	6	ARG	2.7
1	G	161	HIS	2.7
1	E	203	GLU	2.7
1	G	50	ASP	2.6
1	H	47	LEU	2.6
1	H	192	LEU	2.6
1	H	191	SER	2.6
1	G	256	GLU	2.5
1	G	112	PHE	2.5
1	G	162	GLY	2.4
1	B	159	VAL	2.4
1	B	160	TYR	2.3
1	D	99	MET	2.3
1	H	51	VAL	2.3
1	B	255	ILE	2.3
1	E	175	ASN	2.3
1	E	256	GLU	2.2
1	D	256	GLU	2.2
1	G	195	ARG	2.2
1	H	157	THR	2.2
1	E	184	LEU	2.2
1	C	255	ILE	2.1
1	E	160	TYR	2.1
1	F	255	ILE	2.1
1	H	240	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	G	10	GLU	2.0
1	F	175	ASN	2.0
1	H	45	LEU	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	SO4	F	308	5/5	0.98	0.12	0.34	39,41,45,45	0
2	SO4	D	306	5/5	0.99	0.10	-0.20	23,27,29,32	0
2	SO4	G	309	5/5	0.95	0.14	-0.46	61,61,61,62	0
2	SO4	C	303	5/5	0.99	0.07	-1.11	33,34,37,37	0
2	SO4	C	304	5/5	0.98	0.10	-	45,46,46,48	0
2	SO4	B	302	5/5	0.98	0.08	-	42,44,46,47	0
2	SO4	H	310	5/5	0.98	0.09	-	45,46,46,48	0
2	SO4	A	301	5/5	0.98	0.16	-	20,20,20,20	0
2	SO4	F	307	5/5	0.97	0.11	-	52,54,54,55	0
2	SO4	D	305	5/5	0.99	0.07	-	37,41,42,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.