



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:58 PM GMT

PDB ID : 5ELS
Title : Structure of the KH domain of T-STAR in complex with AAAUAA RNA
Authors : Dominguez, C.; Feracci, M.
Deposited on : 2015-11-05
Resolution : 2.87 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

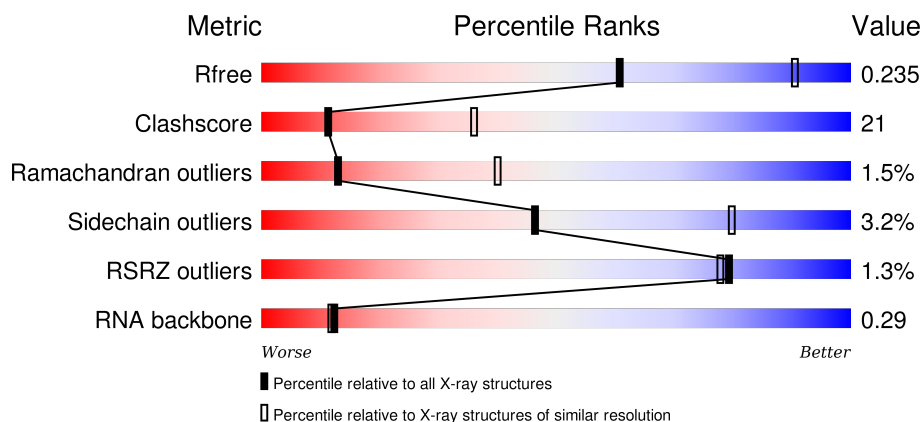
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 2.87 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	1945 (2.90-2.86)
Clashscore	102246	2202 (2.90-2.86)
Ramachandran outliers	100387	2149 (2.90-2.86)
Sidechain outliers	100360	2152 (2.90-2.86)
RSRZ outliers	91569	1950 (2.90-2.86)
RNA backbone	2183	1033 (3.26-2.50)

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	113	 69% 29% .
1	B	113	 67% 27% 5%
1	C	113	 63% 35% .
1	D	113	 58% 36% . . .

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Mol	Chain	Length	Quality of chain
1	E	113	<div><div>%</div><div><div></div><div>53%</div><div>42%</div><div></div><div>.</div><div>.</div></div></div>
1	F	113	<div><div>7%</div><div><div></div><div>43%</div><div>45%</div><div>10%</div><div>.</div><div>.</div></div></div>
2	H	6	<div><div></div><div><div></div><div>33%</div><div>17%</div><div>50%</div></div></div>
2	I	6	<div><div></div><div><div></div><div>50%</div><div>33%</div><div>17%</div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5631 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 KH domain-containing, RNA-binding, signal transduction-associated protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	113	Total	C	N	O	S	0	0	0
			902	581	158	159	4			
1	B	113	Total	C	N	O	S	0	0	0
			902	581	158	159	4			
1	C	113	Total	C	N	O	S	0	0	0
			902	581	158	159	4			
1	D	111	Total	C	N	O	S	0	0	0
			882	568	155	155	4			
1	E	110	Total	C	N	O	S	0	0	0
			874	564	154	152	4			
1	F	112	Total	C	N	O	S	0	0	0
			894	577	156	157	4			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	48	GLY	-	expression tag	UNP O75525
A	49	ALA	-	expression tag	UNP O75525
B	48	GLY	-	expression tag	UNP O75525
B	49	ALA	-	expression tag	UNP O75525
C	48	GLY	-	expression tag	UNP O75525
C	49	ALA	-	expression tag	UNP O75525
D	48	GLY	-	expression tag	UNP O75525
D	49	ALA	-	expression tag	UNP O75525
E	48	GLY	-	expression tag	UNP O75525
E	49	ALA	-	expression tag	UNP O75525
F	48	GLY	-	expression tag	UNP O75525
F	49	ALA	-	expression tag	UNP O75525

- Molecule 2 is a RNA chain called RNA (5'-R(P*AP*AP*AP*UP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	6	Total	C	N	O	P	0	6	0
			130	59	27	38	6			
2	I	6	Total	C	N	O	P	0	6	0
			130	59	27	38	6			

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

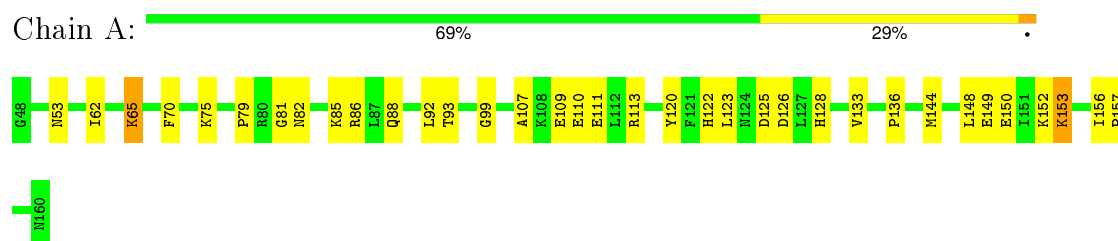


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

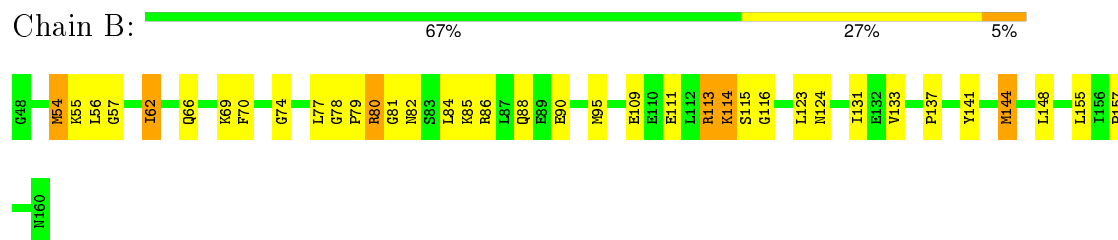
3 Residue-property plots

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.

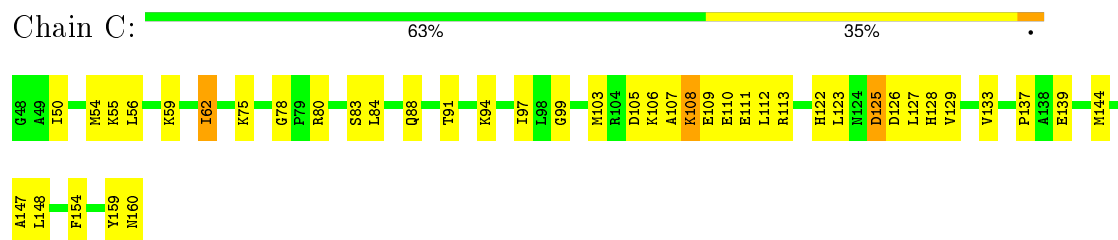
- Molecule 1: KH domain-containing, RNA-binding, signal transduction-associated protein 3



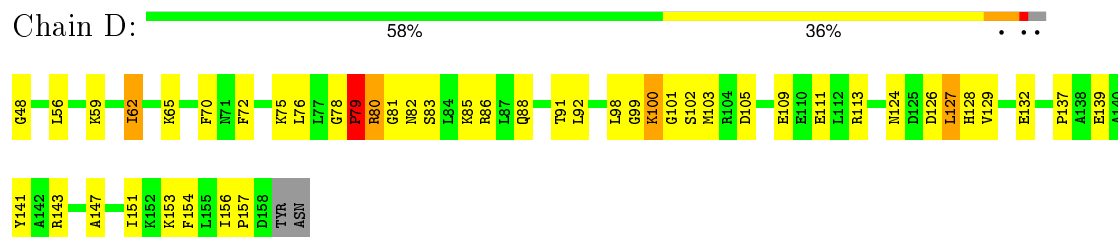
- Molecule 1: KH domain-containing, RNA-binding, signal transduction-associated protein 3



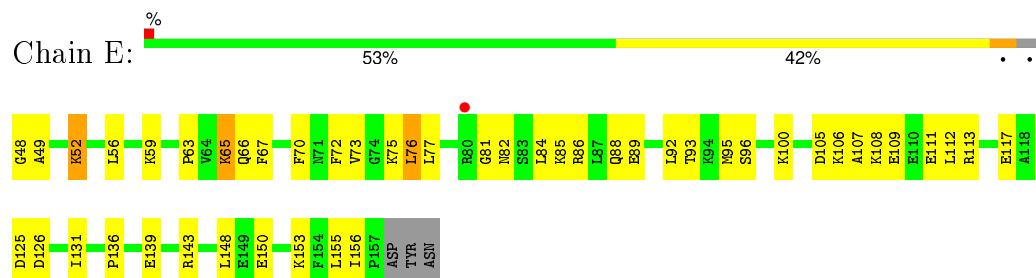
- Molecule 1: KH domain-containing, RNA-binding, signal transduction-associated protein 3



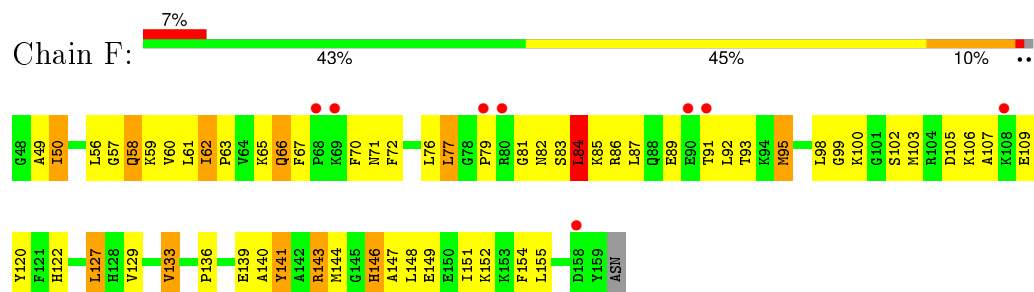
- Molecule 1: KH domain-containing, RNA-binding, signal transduction-associated protein 3



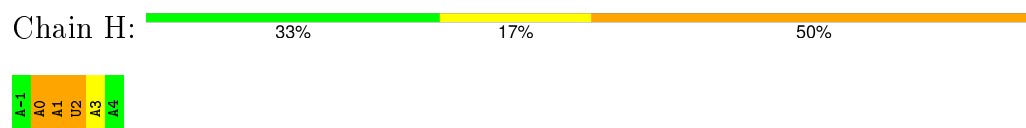
- Molecule 1: KH domain-containing, RNA-binding, signal transduction-associated protein 3



- Molecule 1: KH domain-containing, RNA-binding, signal transduction-associated protein 3



- Molecule 2: RNA (5'-R(P*AP*AP*AP*UP*AP*A)-3')



- Molecule 2: RNA (5'-R(P*AP*AP*AP*UP*AP*A)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	93.73Å 162.22Å 113.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.87 – 2.87 46.87 – 2.87	Depositor EDS
% Data completeness (in resolution range)	99.8 (46.87-2.87) 97.8 (46.87-2.87)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.17	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 2.86Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.188 , 0.236 0.186 , 0.235	Depositor DCC
R_{free} test set	1005 reflections (5.12%)	DCC
Wilson B-factor (Å ²)	43.1	Xtriage
Anisotropy	0.467	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 15.6	EDS
Estimated twinning fraction	0.250 for 1/2*h+1/2*k,3/2*h-1/2*k,-l 0.218 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.217 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
Reported twinning fraction	0.250 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Depositor
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtriage
Outliers	0 of 19995 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5631	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.59% 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	0.64	1/919 (0.1%)	0.76	1/1228 (0.1%)
1	B	0.60	1/919 (0.1%)	0.79	1/1228 (0.1%)
1	C	0.69	1/919 (0.1%)	0.76	1/1228 (0.1%)
1	D	0.59	1/898 (0.1%)	0.80	1/1199 (0.1%)
1	E	0.67	2/890 (0.2%)	0.77	2/1188 (0.2%)
1	F	0.75	5/911 (0.5%)	0.99	8/1217 (0.7%)
2	H	0.43	0/146	0.97	0/225
2	I	0.35	0/146	0.74	0/225
All	All	0.65	11/5748 (0.2%)	0.82	14/7738 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1
1	F	0	2
All	All	0	3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	62	ILE	C-N	11.92	1.56	1.34
1	A	62	ILE	C-N	9.76	1.52	1.34
1	E	113	ARG	CZ-NH1	-8.71	1.21	1.33
1	E	113	ARG	NE-CZ	-8.30	1.22	1.33
1	D	62	ILE	C-N	7.71	1.49	1.34
1	B	62	ILE	C-N	7.56	1.48	1.34
1	F	62	ILE	C-N	-7.17	1.20	1.34
1	F	141	TYR	CD1-CE1	-7.17	1.28	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	141	TYR	CE2-CZ	5.36	1.45	1.38
1	F	141	TYR	CD2-CE2	-5.23	1.31	1.39
1	F	66	GLN	CD-OE1	-5.05	1.12	1.24

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	113	ARG	NE-CZ-NH1	7.96	124.28	120.30
1	F	84	LEU	CB-CG-CD1	7.51	123.77	111.00
1	F	133	VAL	CG1-CB-CG2	-7.49	98.91	110.90
1	F	141	TYR	CA-CB-CG	-7.20	99.72	113.40
1	A	153	LYS	CD-CE-NZ	-6.79	96.09	111.70
1	F	77	LEU	CA-CB-CG	6.66	130.61	115.30
1	F	84	LEU	CB-CG-CD2	-6.18	100.49	111.00
1	E	113	ARG	NE-CZ-NH1	-6.16	117.22	120.30
1	D	127	LEU	CA-CB-CG	6.02	129.15	115.30
1	F	84	LEU	CA-CB-CG	5.89	128.85	115.30
1	C	125	ASP	CB-CG-OD1	5.75	123.48	118.30
1	E	76	LEU	CA-CB-CG	5.44	127.81	115.30
1	F	143	ARG	NE-CZ-NH1	5.02	122.81	120.30
1	F	58	GLN	CA-CB-CG	-5.01	102.37	113.40

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	79	PRO	Peptide
1	F	127	LEU	Peptide
1	F	84	LEU	Peptide

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	902	0	941	20	0
1	B	902	0	941	30	0
1	C	902	0	941	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	882	0	926	47	0
1	E	874	0	922	45	0
1	F	894	0	935	63	0
2	H	130	0	58	4	0
2	I	130	0	57	3	0
3	A	5	0	0	0	0
3	B	10	0	0	1	0
All	All	5631	0	5721	234	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 21.

All (234) 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:F:77:LEU:HD13	1:F:84:LEU:HD23	1.20	1.16
1:F:77:LEU:CD1	1:F:84:LEU:HD23	1.90	1.01
1:A:150:GLU:O	1:A:153:LYS:NZ	1.97	0.95
1:E:84:LEU:HD22	1:E:88:GLN:HE21	1.30	0.94
1:B:86:ARG:NH1	3:B:202:SO4:O4	2.06	0.89
1:B:77:LEU:HB3	1:B:81:GLY:HA2	1.52	0.89
1:F:56:LEU:HB2	1:F:133:VAL:HG12	1.53	0.88
1:D:100:LYS:O	1:D:102:SER:N	2.09	0.86
1:F:83:SER:HA	1:F:86:ARG:HB2	1.57	0.86
1:D:80:ARG:HG3	1:D:82:ASN:H	1.39	0.85
1:D:113:ARG:NH1	1:D:124:ASN:OD1	2.09	0.85
1:F:67:PHE:HB3	1:F:70:PHE:HB3	1.57	0.85
1:F:103:MET:HB2	1:F:109:GLU:HB2	1.61	0.82
1:D:80:ARG:HH12	1:D:86:ARG:HH21	1.27	0.80
1:D:48:GLY:N	1:E:59:LYS:O	2.15	0.79
1:F:77:LEU:HA	1:F:84:LEU:HD22	1.66	0.78
1:D:100:LYS:HD2	1:D:127:LEU:HB3	1.66	0.77
1:C:80:ARG:HA	2:I:3[B]:U:H5"	1.65	0.77
1:F:77:LEU:HA	1:F:84:LEU:CD2	2.14	0.76
1:F:149:GLU:OE2	1:F:152:LYS:NZ	2.19	0.76
1:F:86:ARG:HA	1:F:89:GLU:OE1	1.87	0.75
1:F:92:LEU:HB2	1:F:143:ARG:HD3	1.69	0.75
1:A:149:GLU:OE2	1:A:152:LYS:NZ	2.21	0.74
1:B:70:PHE:HB2	1:C:159:TYR:HB2	1.70	0.74
1:D:100:LYS:HE3	1:D:126:ASP:HB3	1.68	0.73
1:F:139:GLU:O	1:F:143:ARG:HG3	1.89	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:150:GLU:OE1	1:E:153:LYS:NZ	2.22	0.72
1:E:70:PHE:HZ	1:E:75:LYS:HG3	1.55	0.72
1:F:56:LEU:HB2	1:F:133:VAL:CG1	2.20	0.71
1:F:82:ASN:HA	1:F:85:LYS:HB3	1.72	0.71
1:F:95:MET:SD	1:F:95:MET:N	2.61	0.71
1:F:56:LEU:HD21	1:F:141:TYR:CE1	2.25	0.71
1:D:82:ASN:O	1:D:85:LYS:HG2	1.91	0.70
1:F:82:ASN:HB3	1:F:85:LYS:HD3	1.74	0.69
1:F:93:THR:HG21	1:F:147:ALA:HB2	1.73	0.69
1:C:110:GLU:OE2	1:C:113:ARG:NH1	2.24	0.69
1:C:105:ASP:OD2	1:C:106:LYS:N	2.24	0.69
1:A:110:GLU:OE2	1:A:113:ARG:NH1	2.25	0.69
1:F:66:GLN:H	1:F:66:GLN:CD	1.95	0.68
1:E:84:LEU:HD21	1:E:95:MET:HB2	1.76	0.68
1:F:63:PRO:HD2	1:F:155:LEU:HD12	1.74	0.68
1:D:79:PRO:O	1:D:81:GLY:N	2.26	0.67
1:E:82:ASN:HA	1:E:85:LYS:HB3	1.76	0.67
1:C:103:MET:HE1	1:C:112:LEU:HD12	1.76	0.67
1:F:77:LEU:HG	1:F:81:GLY:HA2	1.75	0.67
1:A:109:GLU:HG3	1:A:123:LEU:HD22	1.77	0.66
1:B:114:LYS:O	1:B:116:GLY:N	2.26	0.66
1:A:53:ASN:ND2	1:A:136:PRO:HD3	2.11	0.66
1:F:144:MET:O	1:F:148:LEU:HD12	1.96	0.66
1:F:98:LEU:HD12	1:F:122:HIS:CE1	2.32	0.65
1:E:73:VAL:HA	1:E:76:LEU:HD12	1.78	0.65
1:D:91:THR:HG21	1:D:147:ALA:HA	1.79	0.64
1:C:80:ARG:HA	2:H:3[A]:A:H5"	1.80	0.64
1:D:72:PHE:O	1:D:76:LEU:HD12	1.98	0.64
1:E:105:ASP:OD1	1:E:106:LYS:N	2.31	0.63
1:D:79:PRO:C	1:D:80:ARG:HG2	2.20	0.62
1:C:125:ASP:OD1	1:C:126:ASP:N	2.33	0.62
1:D:62:ILE:O	1:D:100:LYS:NZ	2.33	0.62
1:A:120:TYR:HB3	1:A:123:LEU:HD12	1.82	0.62
1:B:84:LEU:O	1:B:88:GLN:HG3	2.00	0.62
1:F:63:PRO:CD	1:F:155:LEU:HD12	2.31	0.61
1:E:66:GLN:NE2	1:E:67:PHE:CD2	2.68	0.61
1:E:105:ASP:OD1	1:E:107:ALA:N	2.32	0.61
1:D:62:ILE:HD12	1:D:127:LEU:HG	1.81	0.61
1:F:112:LEU:HD23	1:F:120:TYR:CZ	2.35	0.61
1:C:84:LEU:O	1:C:88:GLN:HG3	2.00	0.61
1:B:131:ILE:HD12	1:B:148:LEU:HD23	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:62:ILE:HG21	1:F:127:LEU:HD23	1.84	0.60
1:E:77:LEU:HB3	1:E:81:GLY:HA2	1.83	0.59
1:C:109:GLU:HG3	1:C:123:LEU:HD22	1.84	0.59
1:B:79:PRO:HB2	1:B:80:ARG:HD2	1.85	0.58
1:D:98:LEU:O	1:D:127:LEU:HD12	2.03	0.58
1:C:139:GLU:OE1	1:C:139:GLU:N	2.36	0.58
1:B:79:PRO:C	1:B:80:ARG:HD2	2.24	0.57
1:C:54:MET:HG2	1:C:55:LYS:N	2.19	0.57
1:B:70:PHE:CE1	1:B:157:PRO:HA	2.40	0.57
1:F:89:GLU:OE2	1:F:89:GLU:N	2.37	0.56
1:D:88:GLN:O	1:D:92:LEU:N	2.38	0.56
1:F:106:LYS:NZ	1:F:109:GLU:OE1	2.37	0.56
1:E:108:LYS:O	1:E:111:GLU:N	2.38	0.56
1:D:98:LEU:HD12	1:D:128:HIS:NE2	2.20	0.56
1:E:59:LYS:NZ	1:E:125:ASP:OD1	2.31	0.56
1:E:100:LYS:HE3	1:E:126:ASP:HA	1.86	0.56
1:C:78:GLY:O	2:H:2[A]:U:O2'	2.23	0.56
1:B:66:GLN:N	1:B:66:GLN:OE1	2.38	0.56
1:D:80:ARG:NH1	1:D:86:ARG:HH21	2.01	0.56
1:F:72:PHE:O	1:F:76:LEU:HD12	2.06	0.56
1:D:103:MET:HB2	1:D:109:GLU:HB2	1.88	0.56
1:D:78:GLY:HA3	1:D:80:ARG:HD2	1.86	0.56
1:C:144:MET:O	1:C:148:LEU:HD12	2.06	0.56
1:C:103:MET:HE2	1:C:109:GLU:HB2	1.87	0.55
1:F:70:PHE:HE1	1:F:72:PHE:CD1	2.25	0.55
1:F:91:THR:HG21	1:F:147:ALA:HA	1.87	0.55
1:E:82:ASN:O	1:E:86:ARG:HG2	2.07	0.55
1:E:88:GLN:HB3	1:E:93:THR:O	2.06	0.55
1:F:76:LEU:HG	1:F:154:PHE:HB3	1.89	0.54
1:E:89:GLU:N	1:E:89:GLU:OE1	2.30	0.54
1:E:131:ILE:HD12	1:E:148:LEU:HD23	1.88	0.54
1:D:98:LEU:HB2	1:D:128:HIS:CE1	2.42	0.54
1:A:53:ASN:HD22	1:A:136:PRO:HD3	1.71	0.54
1:E:73:VAL:O	1:E:77:LEU:HD12	2.08	0.54
1:F:99:GLY:HA3	1:F:127:LEU:HA	1.91	0.53
1:C:144:MET:HG3	1:C:148:LEU:HD11	1.89	0.53
1:F:49:ALA:O	1:F:50:ILE:HG22	2.08	0.53
1:E:73:VAL:HA	1:E:76:LEU:CD1	2.38	0.53
1:F:105:ASP:O	1:F:109:GLU:N	2.32	0.53
1:D:137:PRO:HB2	1:E:148:LEU:HD13	1.90	0.53
1:A:125:ASP:OD2	1:A:126:ASP:N	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:70:PHE:HE1	1:F:72:PHE:CE1	2.27	0.53
1:B:56:LEU:HB2	1:B:133:VAL:HG12	1.91	0.53
1:F:100:LYS:O	1:F:127:LEU:HD12	2.09	0.53
1:C:50:ILE:O	1:C:137:PRO:HG2	2.09	0.52
1:E:63:PRO:HB2	1:E:66:GLN:HG3	1.91	0.52
1:F:89:GLU:CD	1:F:89:GLU:H	2.10	0.52
1:E:84:LEU:CD2	1:E:88:GLN:HE21	2.13	0.52
1:D:100:LYS:HD2	1:D:127:LEU:CB	2.36	0.52
1:D:82:ASN:HB3	1:D:85:LYS:HZ2	1.75	0.52
1:A:75:LYS:NZ	1:A:156:ILE:O	2.43	0.51
1:B:74:GLY:O	1:B:78:GLY:N	2.44	0.51
1:C:105:ASP:CG	1:C:107:ALA:H	2.13	0.51
1:D:98:LEU:O	1:D:127:LEU:HA	2.11	0.51
1:D:70:PHE:CD2	1:D:72:PHE:CE2	2.98	0.50
1:B:109:GLU:HG3	1:B:123:LEU:HD13	1.94	0.50
1:A:65:LYS:H	1:A:65:LYS:HD2	1.76	0.50
1:D:129:VAL:HG21	1:D:151:ILE:HD11	1.94	0.50
1:D:100:LYS:HG3	1:D:127:LEU:H	1.77	0.50
1:E:139:GLU:O	1:E:143:ARG:HG3	2.12	0.50
1:D:80:ARG:CD	1:D:83:SER:HB3	2.42	0.50
1:A:88:GLN:HB3	1:A:93:THR:O	2.12	0.50
1:F:110:GLU:O	1:F:113:ARG:N	2.45	0.50
1:C:59:LYS:HG2	1:C:128:HIS:CD2	2.46	0.50
1:B:79:PRO:O	1:B:81:GLY:N	2.45	0.49
1:D:80:ARG:HH12	1:D:86:ARG:NH2	2.04	0.49
1:C:75:LYS:HE2	1:C:154:PHE:HA	1.94	0.49
1:F:98:LEU:HB3	1:F:122:HIS:HB2	1.95	0.49
1:F:63:PRO:C	1:F:66:GLN:OE1	2.51	0.49
1:C:91:THR:HG21	1:C:147:ALA:HA	1.95	0.49
1:D:85:LYS:HZ1	1:D:86:ARG:CZ	2.26	0.48
1:F:56:LEU:HD13	1:F:140:ALA:HB1	1.96	0.48
1:D:139:GLU:O	1:D:143:ARG:HG3	2.12	0.48
1:E:49:ALA:O	1:E:52:LYS:HG3	2.13	0.48
1:E:117:GLU:OE2	1:E:119:LYS:HE2	2.13	0.48
1:D:99:GLY:HA3	1:D:127:LEU:HA	1.94	0.48
1:F:112:LEU:HD23	1:F:120:TYR:CE1	2.48	0.48
1:B:62:ILE:HD13	1:B:155:LEU:HD22	1.95	0.48
1:A:99:GLY:HA2	1:A:122:HIS:O	2.14	0.48
1:A:82:ASN:O	1:A:86:ARG:HG3	2.14	0.48
1:F:136:PRO:HB2	1:F:139:GLU:OE1	2.13	0.47
1:B:54:MET:CE	1:B:137:PRO:HA	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:80:ARG:HD3	1:D:83:SER:N	2.28	0.47
1:F:91:THR:HG22	1:F:146:HIS:ND1	2.29	0.47
1:F:60:VAL:HG12	1:F:129:VAL:O	2.14	0.47
1:F:105:ASP:OD2	1:F:107:ALA:HB3	2.14	0.47
1:F:144:MET:C	1:F:148:LEU:HD12	2.34	0.47
1:B:57:GLY:HA2	1:B:131:ILE:O	2.14	0.47
1:C:88:GLN:HE21	1:C:94:LYS:HA	1.80	0.46
1:A:93:THR:HG22	1:A:133:VAL:HG22	1.97	0.46
1:D:80:ARG:O	1:D:82:ASN:ND2	2.49	0.46
1:D:80:ARG:HD2	1:D:83:SER:HB3	1.98	0.46
1:F:57:GLY:C	1:F:58:GLN:HG2	2.36	0.46
2:H:0[A]:A:H4'	2:H:1[A]:A:O5'	2.14	0.46
1:C:105:ASP:HB3	1:C:108:LYS:HD2	1.97	0.46
1:F:77:LEU:CA	1:F:84:LEU:CD2	2.90	0.46
1:B:88:GLN:HG2	1:B:95:MET:SD	2.56	0.46
1:F:144:MET:HG2	1:F:148:LEU:HD11	1.98	0.45
1:C:59:LYS:NZ	1:C:125:ASP:OD2	2.33	0.45
1:C:56:LEU:HB2	1:C:133:VAL:HG12	1.97	0.45
1:D:72:PHE:O	1:D:75:LYS:N	2.49	0.45
1:D:80:ARG:HG3	1:D:82:ASN:N	2.20	0.45
1:B:70:PHE:HB2	1:C:159:TYR:CB	2.45	0.45
1:E:65:LYS:HD2	1:E:65:LYS:C	2.36	0.45
1:C:62:ILE:HD12	1:C:127:LEU:HG	1.98	0.45
1:B:69:LYS:O	1:C:160:ASN:HB3	2.17	0.45
1:C:78:GLY:HA2	2:I:2[B]:A:N3	2.31	0.45
1:E:155:LEU:O	1:E:156:ILE:HG13	2.16	0.45
1:E:85:LYS:O	1:E:89:GLU:OE1	2.34	0.45
1:B:114:LYS:O	1:B:114:LYS:HD2	2.16	0.45
1:D:70:PHE:CD2	1:D:72:PHE:HE2	2.35	0.45
1:E:49:ALA:HB3	1:E:52:LYS:CG	2.47	0.45
1:F:57:GLY:O	1:F:58:GLN:HG2	2.15	0.45
1:F:143:ARG:O	1:F:146:HIS:HB3	2.16	0.45
1:F:110:GLU:HG3	1:F:113:ARG:NH1	2.31	0.45
1:B:86:ARG:O	1:B:90:GLU:HG3	2.17	0.45
1:E:108:LYS:HG3	1:E:109:GLU:N	2.32	0.45
1:E:66:GLN:NE2	1:E:67:PHE:CG	2.85	0.44
1:A:144:MET:O	1:A:148:LEU:HG	2.17	0.44
1:E:136:PRO:HB2	1:E:139:GLU:OE1	2.17	0.44
1:B:80:ARG:HB3	1:B:82:ASN:ND2	2.32	0.44
1:D:85:LYS:NZ	1:D:86:ARG:CZ	2.80	0.44
1:E:72:PHE:C	1:E:76:LEU:HD12	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:ASP:O	1:A:128:HIS:ND1	2.36	0.44
1:D:100:LYS:HE3	1:D:127:LEU:H	1.82	0.43
1:D:141:TYR:HE2	1:E:56:LEU:HD13	1.83	0.43
1:A:107:ALA:O	1:A:111:GLU:HG3	2.18	0.43
1:C:78:GLY:HA2	2:I:2[B]:A:C2	2.54	0.43
1:F:84:LEU:N	1:F:84:LEU:HD22	2.34	0.43
1:F:62:ILE:CG2	1:F:127:LEU:HD23	2.48	0.43
1:E:65:LYS:HE3	1:E:66:GLN:HB3	2.00	0.43
1:F:62:ILE:HA	1:F:63:PRO:HD3	1.79	0.43
1:C:107:ALA:O	1:C:111:GLU:HG3	2.18	0.43
1:A:70:PHE:CE1	1:A:157:PRO:HA	2.53	0.43
1:D:82:ASN:CB	1:D:85:LYS:HZ2	2.31	0.43
1:F:106:LYS:HA	1:F:109:GLU:HB3	2.01	0.43
1:F:111:GLU:HG2	1:F:112:LEU:HD12	2.00	0.42
1:A:88:GLN:O	1:A:92:LEU:N	2.52	0.42
1:F:93:THR:HG21	1:F:147:ALA:CB	2.45	0.42
1:B:113:ARG:NH2	1:B:124:ASN:OD1	2.48	0.42
1:E:120:TYR:HB3	1:E:123:LEU:HD12	2.01	0.42
1:B:82:ASN:HA	1:B:85:LYS:HG2	2.01	0.42
1:F:148:LEU:HA	1:F:151:ILE:HG22	2.02	0.42
1:D:59:LYS:O	1:E:48:GLY:N	2.53	0.42
1:E:82:ASN:HA	1:E:85:LYS:CB	2.48	0.41
1:D:153:LYS:O	1:D:156:ILE:HG13	2.20	0.41
1:E:70:PHE:CZ	1:E:75:LYS:HG3	2.44	0.41
1:B:79:PRO:CB	1:B:80:ARG:HD2	2.50	0.41
1:C:78:GLY:HA2	2:H:2[A]:U:C6	2.55	0.41
1:B:144:MET:O	1:B:148:LEU:HG	2.20	0.41
1:E:109:GLU:HA	1:E:112:LEU:HD13	2.03	0.41
1:C:99:GLY:HA2	1:C:122:HIS:O	2.21	0.41
1:F:61:LEU:HD12	1:F:62:ILE:H	1.86	0.41
1:B:54:MET:HG2	1:B:55:LYS:N	2.35	0.41
1:D:153:LYS:HG3	1:D:154:PHE:CD1	2.56	0.41
1:E:67:PHE:HD2	1:E:72:PHE:HZ	1.68	0.41
1:E:139:GLU:OE1	1:E:139:GLU:N	2.52	0.41
1:F:87:LEU:HA	1:F:87:LEU:HD23	1.66	0.41
1:C:144:MET:C	1:C:148:LEU:HD12	2.42	0.41
1:C:97:ILE:HG12	1:C:129:VAL:HG22	2.03	0.41
1:E:88:GLN:O	1:E:92:LEU:N	2.54	0.40
1:B:79:PRO:HB2	1:B:80:ARG:CD	2.50	0.40
1:E:65:LYS:CE	1:E:66:GLN:HB3	2.51	0.40
1:C:78:GLY:HA3	1:C:83:SER:OG	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:PRO:O	1:A:81:GLY:N	2.54	0.40
1:B:56:LEU:HD11	1:B:141:TYR:CD1	2.57	0.40
1:D:56:LEU:O	1:D:132:GLU:HA	2.21	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	111/113 (98%)	108 (97%)	3 (3%)	0	100	100
1	B	111/113 (98%)	107 (96%)	2 (2%)	2 (2%)	11	35
1	C	111/113 (98%)	108 (97%)	3 (3%)	0	100	100
1	D	109/113 (96%)	102 (94%)	1 (1%)	6 (6%)	2	6
1	E	108/113 (96%)	104 (96%)	4 (4%)	0	100	100
1	F	110/113 (97%)	98 (89%)	10 (9%)	2 (2%)	11	35
All	All	660/678 (97%)	627 (95%)	23 (4%)	10 (2%)	13	40

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	115	SER
1	D	79	PRO
1	D	80	ARG
1	D	101	GLY
1	D	157	PRO
1	F	79	PRO
1	F	50	ILE
1	B	80	ARG
1	D	100	LYS
1	D	105	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	96/96 (100%)	94 (98%)	2 (2%)	61	87
1	B	96/96 (100%)	92 (96%)	4 (4%)	36	71
1	C	96/96 (100%)	95 (99%)	1 (1%)	82	95
1	D	94/96 (98%)	92 (98%)	2 (2%)	61	87
1	E	93/96 (97%)	90 (97%)	3 (3%)	46	80
1	F	95/96 (99%)	89 (94%)	6 (6%)	22	52
All	All	570/576 (99%)	552 (97%)	18 (3%)	46	80

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	LYS
1	A	85	LYS
1	B	54	MET
1	B	111	GLU
1	B	114	LYS
1	B	144	MET
1	C	108	LYS
1	D	65	LYS
1	D	111	GLU
1	E	52	LYS
1	E	65	LYS
1	E	96	SER
1	F	59	LYS
1	F	65	LYS
1	F	71	ASN
1	F	95	MET
1	F	102	SER
1	F	146	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	ASN
1	A	71	ASN
1	A	124	ASN
1	B	58	GLN
1	C	53	ASN
1	C	88	GLN
1	D	82	ASN
1	E	82	ASN
1	E	88	GLN
1	F	53	ASN
1	F	71	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	H	5/6 (83%)	3 (60%)	0
2	I	5/6 (83%)	2 (40%)	0
All	All	10/12 (83%)	5 (50%)	0

All (5) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	H	0[A]	A
2	H	1[A]	A
2	H	2[A]	U
2	I	1[B]	A
2	I	2[B]	A

There are no RNA pucker outliers to report.

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

3 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	201	-	4,4,4	0.20	0	6,6,6	0.24	0
3	SO4	B	201	-	4,4,4	0.27	0	6,6,6	0.32	0
3	SO4	B	202	-	4,4,4	0.41	0	6,6,6	0.35	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	201	-	-	0/0/0/0	0/0/0/0
3	SO4	B	201	-	-	0/0/0/0	0/0/0/0
3	SO4	B	202	-	-	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	202	SO4	1	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 ⓘ

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	113/113 (100%)	-0.48	0 100 100	17, 28, 46, 53	0
1	B	113/113 (100%)	-0.31	0 100 100	23, 39, 80, 95	0
1	C	113/113 (100%)	-0.47	0 100 100	14, 27, 69, 82	0
1	D	111/113 (98%)	-0.20	0 100 100	25, 51, 88, 113	0
1	E	110/113 (97%)	-0.18	1 (0%) 85 84	35, 57, 104, 133	0
1	F	112/113 (99%)	0.38	8 (7%) 19 13	47, 82, 132, 189	0
2	H	6/6 (100%)	0.02	0 100 100	31, 39, 47, 47	6 (100%)
2	I	6/6 (100%)	-0.06	0 100 100	30, 38, 44, 47	6 (100%)
All	All	684/690 (99%)	-0.21	9 (1%) 79 78	14, 46, 106, 189	12 (1%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	158	ASP	3.4
1	F	80	ARG	3.4
1	F	90	GLU	3.3
1	F	108	LYS	3.0
1	F	69	LYS	2.7
1	F	68	PRO	2.7
1	F	91	THR	2.5
1	E	80	ARG	2.4
1	F	79	PRO	2.2

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

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(\AA^2)	Q<0.9
3	SO4	B	202	5/5	0.96	0.17	-1.07	44,45,48,60	0
3	SO4	B	201	5/5	0.97	0.14	-	54,56,61,73	0
3	SO4	A	201	5/5	0.96	0.14	-	59,59,67,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.