



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:17 AM GMT

PDB ID : 2WJV
Title : CRYSTAL STRUCTURE OF THE COMPLEX BETWEEN HUMAN NON-SENSE MEDIATED DECAY FACTORS UPF1 AND UPF2
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Deposited on : 2009-06-01
Resolution : 2.85 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

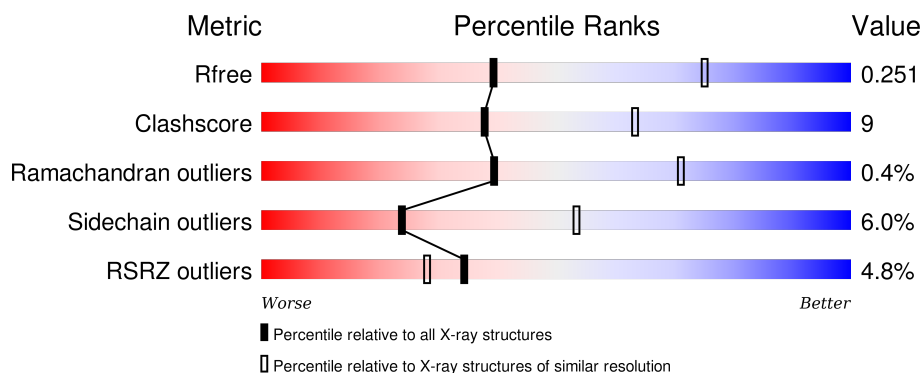
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	2228 (2.90-2.82)
Clashscore	102246	2499 (2.90-2.82)
Ramachandran outliers	100387	2439 (2.90-2.82)
Sidechain outliers	100360	2442 (2.90-2.82)
RSRZ outliers	91569	2236 (2.90-2.82)

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	800	<div> <div>4%</div> <div>78%</div> <div>17%</div> <div>• •</div> </div>
1	B	800	<div> <div>4%</div> <div>77%</div> <div>17%</div> <div>• •</div> </div>
2	D	97	<div> <div>2%</div> <div>47%</div> <div>6%</div> <div>•</div> <div>44%</div> </div>
2	E	97	<div> <div>10%</div> <div>43%</div> <div>15%</div> <div>•</div> <div>38%</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
4	SO4	A	994	-	-	X	-
4	SO4	A	996	-	-	-	X
4	SO4	A	997	-	-	-	X
4	SO4	A	998	-	-	X	-
4	SO4	A	999	-	-	-	X
4	SO4	B	996	-	-	-	X
4	SO4	B	999	-	-	X	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13187 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 REGULATOR OF NONSENSE TRANSCRIPTS 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	772	Total	C	N	O	S	0	0	0
			6094	3867	1067	1125	35			
1	B	768	Total	C	N	O	S	0	0	0
			6073	3858	1065	1115	35			

- Molecule 2 is a protein called REGULATOR OF NONSENSE TRANSCRIPTS 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	54	Total	C	N	O	S	0	0	0
			426	268	70	81	7			
2	E	60	Total	C	N	O	S	0	0	0
			473	295	82	89	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	1102	ALA	-	EXPRESSION TAG	UNP Q9HAU5
D	1103	MET	-	EXPRESSION TAG	UNP Q9HAU5
D	1104	GLY	-	EXPRESSION TAG	UNP Q9HAU5
E	1102	ALA	-	EXPRESSION TAG	UNP Q9HAU5
E	1103	MET	-	EXPRESSION TAG	UNP Q9HAU5
E	1104	GLY	-	EXPRESSION TAG	UNP Q9HAU5

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total	Zn	0	0
			3	3		
3	A	3	Total	Zn	0	0
			3	3		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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

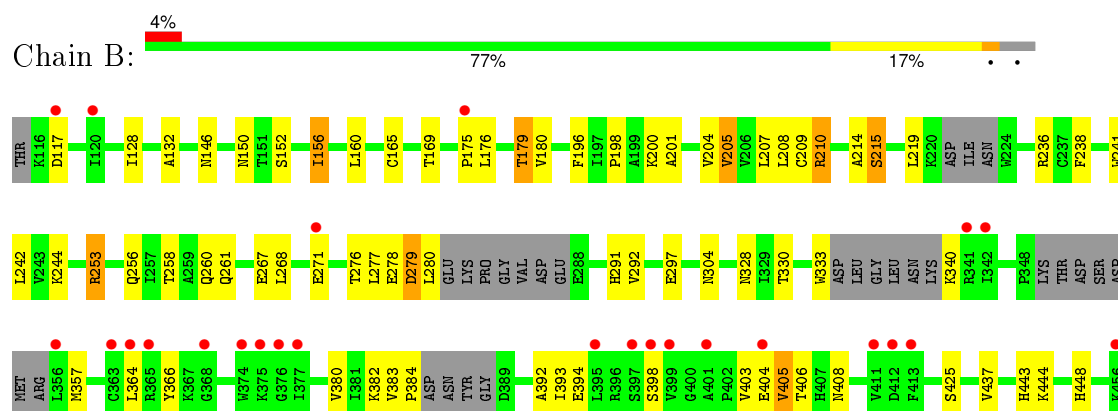
3 Residue-property plots

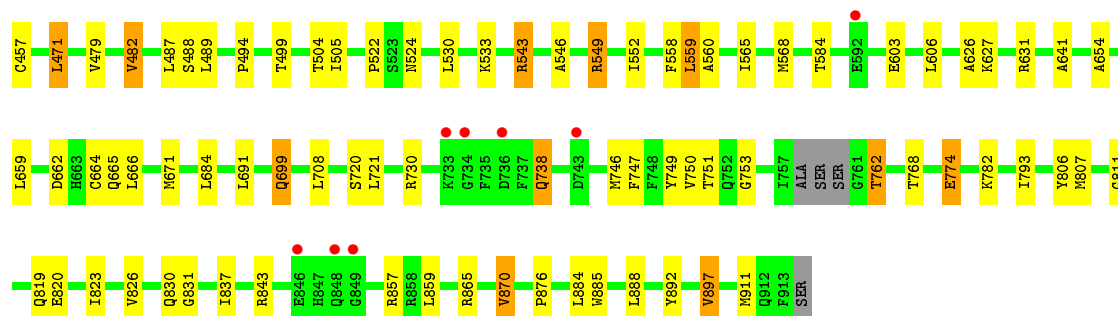
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: REGULATOR OF NONSENSE TRANSCRIPTS 1



• Molecule 1: REGULATOR OF NONSENSE TRANSCRIPTS 1

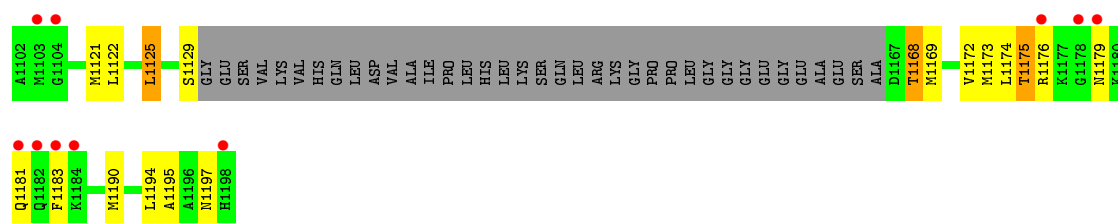
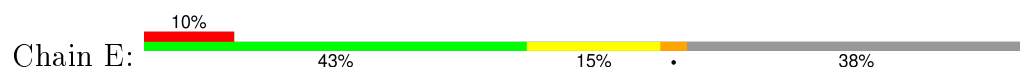




• Molecule 2: REGULATOR OF NONSENSE TRANSCRIPTS 2



• Molecule 2: REGULATOR OF NONSENSE TRANSCRIPTS 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	92.41 Å 97.29 Å 124.60 Å 90.00° 102.38° 90.00°	Depositor
Resolution (Å)	48.79 – 2.85 48.77 – 2.85	Depositor EDS
% Data completeness (in resolution range)	96.9 (48.79-2.85) 96.9 (48.77-2.85)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.86 Å)	Xtriage
Refinement program	REFMAC 5.5.0038	Depositor
R, R_{free}	0.200 , 0.248 0.206 , 0.251	Depositor DCC
R_{free} test set	2463 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	64.9	Xtriage
Anisotropy	0.265	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 48997 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13187	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.78% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 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	0.53	1/6215 (0.0%)	0.61	0/8412
1	B	0.54	1/6194 (0.0%)	0.62	0/8382
2	D	0.45	0/430	0.59	0/576
2	E	0.51	0/478	0.65	0/639
All	All	0.53	2/13317 (0.0%)	0.62	0/18009

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	664	CYS	CB-SG	-5.58	1.72	1.81
1	A	664	CYS	CB-SG	-5.45	1.73	1.81

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6094	0	6140	107	0
1	B	6073	0	6126	119	0
2	D	426	0	421	18	0
2	E	473	0	475	11	0
3	A	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	3	0	0	0	0
4	A	65	0	0	8	0
4	B	50	0	0	4	0
All	All	13187	0	13162	243	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 9.

All (243) 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:206:VAL:CG1	2:D:1174:LEU:HD12	1.94	0.98
1:B:364:LEU:HD22	1:B:393:ILE:HD11	1.46	0.97
1:A:364:LEU:HD22	1:A:393:ILE:HD11	1.49	0.91
1:B:524:ASN:HD22	1:B:543:ARG:HH22	1.15	0.91
1:A:584:THR:HG22	1:A:585:GLY:O	1.71	0.89
1:B:762:THR:O	1:B:762:THR:CG2	2.26	0.84
1:B:268:LEU:HD22	1:B:280:LEU:HD21	1.63	0.81
1:A:524:ASN:HD22	1:A:543:ARG:HH22	1.26	0.80
1:B:196:PHE:CD2	1:B:205:VAL:HG13	2.18	0.78
1:B:762:THR:HG22	1:B:762:THR:O	1.83	0.77
1:A:580:LEU:HD23	1:A:587:LEU:HD22	1.67	0.77
1:A:473:HIS:ND1	4:A:994:SO4:O1	2.16	0.76
2:D:1168:THR:HG23	2:D:1169:MET:N	2.01	0.76
1:A:207:LEU:HD11	1:A:232:LEU:HD13	1.67	0.74
1:A:253:ARG:NH2	1:A:437:VAL:O	2.21	0.74
1:B:382:LYS:HB3	1:B:392:ALA:HB3	1.69	0.73
1:B:443:HIS:HD2	1:B:448:HIS:ND1	1.86	0.73
1:A:382:LYS:HB3	1:A:392:ALA:HB3	1.70	0.73
1:A:482:VAL:HG22	1:A:488:SER:CB	2.19	0.72
1:A:499:THR:HG23	1:A:530:LEU:HD13	1.71	0.72
2:D:1168:THR:CG2	2:D:1169:MET:N	2.55	0.69
1:B:482:VAL:HG22	1:B:488:SER:CB	2.24	0.68
1:B:708:LEU:HD22	1:B:870:VAL:HG13	1.75	0.68
1:A:482:VAL:HG22	1:A:488:SER:HB2	1.75	0.67
1:B:666:LEU:HD12	1:B:830:GLN:CG	2.25	0.67
2:E:1168:THR:CG2	2:E:1169:MET:N	2.56	0.67
1:A:793:ILE:HG21	1:A:826:VAL:HG12	1.77	0.67
1:A:907:ARG:NH2	4:A:988:SO4:O2	2.28	0.66
2:D:1175:THR:O	2:D:1175:THR:HG22	1.95	0.66
1:B:253:ARG:NH2	1:B:437:VAL:O	2.29	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:1168:THR:HG21	2:D:1187:ASN:HB3	1.78	0.66
1:B:499:THR:HG23	1:B:530:LEU:HD13	1.77	0.65
1:B:558:PHE:CE1	1:B:559:LEU:HD13	2.33	0.64
1:A:443:HIS:HD2	1:A:448:HIS:ND1	1.95	0.64
2:D:1124:ASN:HD22	2:D:1124:ASN:N	1.95	0.64
1:A:330:THR:HG23	1:A:406:THR:CG2	2.29	0.63
2:D:1169:MET:HB2	2:D:1190:MET:HE1	1.80	0.63
2:E:1168:THR:HG23	2:E:1169:MET:N	2.14	0.63
1:B:207:LEU:O	1:B:208:LEU:HD23	1.99	0.63
1:B:826:VAL:HG22	4:B:996:SO4:O3	1.97	0.63
1:B:383:VAL:HG12	1:B:384:PRO:HD2	1.81	0.63
1:A:330:THR:HG23	1:A:406:THR:HG23	1.81	0.62
1:A:146:ASN:HD22	1:A:156:ILE:HD11	1.65	0.62
1:B:278:GLU:HG2	1:B:304:ASN:HD21	1.65	0.62
1:B:750:VAL:HG11	1:B:911:MET:HE1	1.82	0.62
1:A:333:TRP:CB	1:A:403:VAL:HG13	2.30	0.61
2:D:1190:MET:HA	2:D:1190:MET:HE2	1.81	0.61
1:A:549:ARG:HA	1:A:552:ILE:HD13	1.82	0.61
1:B:330:THR:HG23	1:B:406:THR:CG2	2.31	0.61
1:B:383:VAL:HG12	1:B:384:PRO:CD	2.31	0.60
1:A:568:MET:HE3	1:A:606:LEU:HD21	1.82	0.60
1:A:383:VAL:HG12	1:A:384:PRO:HD2	1.83	0.60
1:B:699:GLN:HE21	1:B:720:SER:HB2	1.67	0.59
1:B:333:TRP:CB	1:B:403:VAL:HG13	2.33	0.59
1:A:558:PHE:CE1	1:A:559:LEU:HD13	2.37	0.59
1:A:146:ASN:HD22	1:A:156:ILE:CD1	2.14	0.59
1:B:843:ARG:NH1	4:B:991:SO4:O2	2.34	0.59
1:B:210:ARG:O	1:B:214:ALA:HB3	2.03	0.59
1:A:418:THR:HG23	4:A:998:SO4:O3	2.03	0.58
1:B:666:LEU:HD12	1:B:830:GLN:HG2	1.85	0.58
1:B:482:VAL:HG22	1:B:488:SER:HB2	1.86	0.58
1:A:383:VAL:HG12	1:A:384:PRO:CD	2.33	0.58
1:B:793:ILE:HG21	1:B:826:VAL:HG12	1.85	0.58
1:B:549:ARG:HA	1:B:552:ILE:HD13	1.85	0.57
1:A:584:THR:CG2	1:A:585:GLY:N	2.68	0.57
1:A:332:ARG:NH2	1:B:811:GLY:O	2.37	0.57
1:B:330:THR:HG23	1:B:406:THR:HG23	1.86	0.57
1:A:699:GLN:HE21	1:A:720:SER:HB2	1.69	0.57
1:B:146:ASN:HD22	1:B:156:ILE:HD11	1.69	0.57
1:B:626:ALA:O	1:B:627:LYS:HB2	2.05	0.57
1:B:494:PRO:HA	4:B:999:SO4:O2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:666:LEU:HD12	1:B:830:GLN:CD	2.26	0.56
1:A:333:TRP:CG	1:A:403:VAL:HG13	2.41	0.56
1:B:671:MET:HA	1:B:671:MET:CE	2.36	0.56
2:D:1169:MET:HB2	2:D:1190:MET:CE	2.37	0.55
1:B:708:LEU:HD22	1:B:870:VAL:CG1	2.36	0.55
1:A:626:ALA:O	1:A:627:LYS:HB2	2.05	0.55
1:A:897:VAL:HG13	1:A:897:VAL:O	2.06	0.55
1:A:584:THR:HG22	1:A:585:GLY:N	2.22	0.55
1:A:146:ASN:HD21	1:A:238:PHE:H	1.55	0.55
1:A:333:TRP:CE3	1:A:403:VAL:HG22	2.41	0.55
1:A:206:VAL:HG13	2:D:1174:LEU:HD12	1.84	0.54
1:B:888:LEU:HD11	1:B:892:TYR:CE1	2.42	0.54
1:A:671:MET:CE	1:A:671:MET:HA	2.38	0.54
1:B:774:GLU:HG2	1:B:806:TYR:HE2	1.71	0.54
1:B:859:LEU:HD23	1:B:884:LEU:CD2	2.38	0.54
1:B:750:VAL:CG1	1:B:911:MET:HE1	2.37	0.54
1:B:444:LYS:HG2	1:B:487:LEU:HD13	1.89	0.54
1:A:479:VAL:HG22	1:A:505:ILE:HG13	1.89	0.54
1:B:807:MET:HE1	1:B:823:ILE:HD11	1.88	0.54
1:A:774:GLU:HG2	1:A:806:TYR:HE2	1.73	0.53
1:A:206:VAL:HG11	2:D:1174:LEU:HD12	1.86	0.53
1:B:558:PHE:CD1	1:B:559:LEU:HD13	2.44	0.53
1:A:888:LEU:HD11	1:A:892:TYR:CE1	2.44	0.53
1:B:437:VAL:HG12	1:B:654:ALA:HB3	1.91	0.53
1:A:403:VAL:O	1:A:403:VAL:HG12	2.09	0.53
1:A:580:LEU:CD2	1:A:587:LEU:HD22	2.38	0.53
1:A:206:VAL:HG12	2:D:1174:LEU:HD12	1.85	0.52
1:B:196:PHE:HD2	1:B:205:VAL:HG13	1.71	0.51
1:A:546:ALA:HB3	1:A:549:ARG:HG3	1.91	0.51
1:B:366:TYR:OH	1:B:405:VAL:HG11	2.10	0.51
2:D:1168:THR:HG23	2:D:1169:MET:H	1.73	0.51
1:B:179:THR:HG21	2:E:1175:THR:HG21	1.93	0.51
1:B:487:LEU:HD23	1:B:487:LEU:C	2.31	0.51
1:B:831:GLY:HA2	4:B:999:SO4:O3	2.10	0.51
1:B:443:HIS:CD2	1:B:448:HIS:ND1	2.74	0.51
1:A:487:LEU:C	1:A:487:LEU:HD23	2.31	0.51
1:B:333:TRP:CG	1:B:403:VAL:HG13	2.46	0.51
1:B:146:ASN:HD21	1:B:238:PHE:H	1.59	0.51
1:A:799:GLN:O	1:A:803:LEU:HB2	2.10	0.51
1:B:750:VAL:HG11	1:B:911:MET:CE	2.40	0.51
1:B:487:LEU:HD23	1:B:488:SER:N	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:524:ASN:HD22	1:B:543:ARG:NH2	1.97	0.51
1:A:558:PHE:CD1	1:A:559:LEU:HD13	2.46	0.51
1:A:210:ARG:HA	1:A:214:ALA:HB3	1.92	0.51
1:A:640:GLN:NE2	1:A:830:GLN:OE1	2.44	0.50
1:B:546:ALA:HB3	1:B:549:ARG:HG3	1.92	0.50
1:A:666:LEU:HD12	1:A:830:GLN:CG	2.42	0.50
1:B:568:MET:HE3	1:B:606:LEU:HD21	1.91	0.50
2:D:1190:MET:HE2	2:D:1190:MET:CA	2.42	0.50
1:A:366:TYR:OH	1:A:405:VAL:HG11	2.11	0.50
1:B:169:THR:OG1	1:B:180:VAL:HG22	2.12	0.50
1:B:479:VAL:HG22	1:B:505:ILE:HG13	1.93	0.49
1:B:897:VAL:HG13	1:B:897:VAL:O	2.12	0.49
1:A:562:HIS:ND1	4:A:999:SO4:O4	2.38	0.49
1:B:146:ASN:HD22	1:B:156:ILE:CD1	2.26	0.49
1:B:489:LEU:HD22	1:B:659:LEU:HB2	1.94	0.49
1:A:206:VAL:HG12	2:D:1174:LEU:HB2	1.94	0.49
1:A:146:ASN:ND2	1:A:156:ILE:HD11	2.28	0.48
1:A:747:PHE:HA	1:A:897:VAL:HG13	1.95	0.48
1:B:565:ILE:HD11	1:B:603:GLU:HG3	1.94	0.48
1:A:380:VAL:HA	1:A:393:ILE:HG22	1.95	0.48
1:A:328:ASN:ND2	1:A:408:ASN:HD22	2.11	0.48
1:B:156:ILE:HD12	1:B:156:ILE:H	1.79	0.48
1:B:291:HIS:HD2	1:B:292:VAL:O	1.96	0.48
1:A:770:ALA:HB1	1:A:803:LEU:HD13	1.96	0.48
1:B:403:VAL:O	1:B:403:VAL:HG12	2.13	0.48
1:A:152:SER:O	1:A:236:ARG:NH1	2.46	0.48
1:A:903:LEU:HA	1:A:906:LEU:HG	1.96	0.48
1:A:487:LEU:HD23	1:A:488:SER:N	2.29	0.47
2:D:1175:THR:CG2	2:D:1175:THR:O	2.61	0.47
1:A:150:ASN:HD21	1:A:256:GLN:HB3	1.79	0.47
1:B:210:ARG:HG2	2:E:1125:LEU:HD21	1.97	0.47
1:A:708:LEU:HD22	1:A:870:VAL:HG13	1.96	0.47
1:A:437:VAL:HG12	1:A:654:ALA:HB3	1.96	0.47
1:B:747:PHE:HA	1:B:897:VAL:HG13	1.95	0.47
1:A:169:THR:OG1	1:A:180:VAL:HG22	2.15	0.47
1:A:769:GLU:O	1:A:773:VAL:HG23	2.15	0.47
1:B:662:ASP:HB3	1:B:665:GLN:HE21	1.79	0.47
1:B:333:TRP:CE3	1:B:403:VAL:HG22	2.50	0.47
1:A:333:TRP:CD2	1:A:403:VAL:HG22	2.49	0.47
1:B:152:SER:O	1:B:236:ARG:NH1	2.48	0.47
1:B:366:TYR:CZ	1:B:405:VAL:HG11	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:LYS:O	1:B:201:ALA:HB3	2.14	0.46
1:B:297:GLU:HA	1:B:297:GLU:OE1	2.16	0.46
1:A:335:LEU:HD22	1:B:782:LYS:HA	1.97	0.46
1:B:150:ASN:HD21	1:B:256:GLN:HB3	1.81	0.46
1:B:380:VAL:HA	1:B:393:ILE:HG22	1.97	0.46
1:A:156:ILE:H	1:A:156:ILE:HD12	1.80	0.46
1:A:333:TRP:HB2	1:A:403:VAL:HG13	1.98	0.46
1:B:366:TYR:CE1	1:B:405:VAL:HG11	2.51	0.46
1:A:769:GLU:HG3	1:A:873:VAL:HG12	1.98	0.46
1:A:684:LEU:HD22	1:A:688:LEU:CD1	2.46	0.46
1:B:198:PRO:HB3	2:E:1194:LEU:HD11	1.98	0.45
1:A:444:LYS:HG2	1:A:487:LEU:HD13	1.98	0.45
1:B:471:LEU:HD21	1:B:504:THR:HG21	1.98	0.45
1:A:422:ARG:NE	4:A:998:SO4:O3	2.47	0.45
1:A:160:LEU:HD23	1:A:165:CYS:HB2	1.99	0.45
1:A:684:LEU:HD22	1:A:688:LEU:HD11	1.98	0.45
2:E:1176:ARG:HA	2:E:1181:GLN:HA	1.99	0.45
1:B:662:ASP:CB	1:B:665:GLN:HE21	2.29	0.45
1:A:472:ASN:HB2	4:A:994:SO4:O1	2.17	0.45
1:B:176:LEU:HD23	1:B:179:THR:OG1	2.17	0.45
1:A:340:LYS:HD2	1:A:394:GLU:OE2	2.17	0.45
1:B:837:ILE:HG13	1:B:865:ARG:HB2	1.99	0.45
1:B:261:GLN:OE1	1:B:277:LEU:HD21	2.17	0.45
1:B:146:ASN:ND2	1:B:156:ILE:HD11	2.31	0.44
1:B:219:LEU:HD23	1:B:219:LEU:O	2.18	0.44
1:B:543:ARG:O	1:B:560:ALA:HA	2.18	0.44
1:A:403:VAL:CG1	1:A:403:VAL:O	2.66	0.44
2:E:1174:LEU:HD23	2:E:1183:PHE:CD2	2.52	0.44
2:D:1124:ASN:N	2:D:1124:ASN:ND2	2.65	0.44
1:B:738:GLN:HA	1:B:738:GLN:HE21	1.82	0.44
1:A:489:LEU:HD22	1:A:659:LEU:HB2	1.99	0.44
1:A:758:ALA:O	1:A:760:SER:N	2.50	0.44
1:B:328:ASN:ND2	1:B:408:ASN:HD22	2.16	0.44
1:B:132:ALA:HB1	1:B:244:LYS:HG3	1.98	0.44
1:B:205:VAL:HG12	2:E:1173:MET:HG3	2.00	0.44
1:B:333:TRP:HB2	1:B:403:VAL:HG13	1.99	0.43
1:B:522:PRO:HG3	1:B:641:ALA:HB2	1.99	0.43
1:A:366:TYR:CE1	1:A:405:VAL:HG11	2.53	0.43
1:A:366:TYR:CZ	1:A:405:VAL:HG11	2.53	0.43
1:A:471:LEU:HD21	1:A:504:THR:HG21	2.00	0.43
1:A:204:VAL:O	1:A:205:VAL:HG23	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:746:MET:O	1:A:897:VAL:HG21	2.18	0.43
1:B:258:THR:O	1:B:261:GLN:N	2.49	0.43
1:B:160:LEU:HD23	1:B:165:CYS:HB2	1.99	0.43
1:A:837:ILE:HG13	1:A:865:ARG:HB2	1.99	0.43
1:A:196:PHE:CD2	1:A:205:VAL:CG1	3.02	0.43
1:B:888:LEU:CD1	1:B:892:TYR:CE1	3.02	0.43
1:A:826:VAL:HG22	4:A:996:SO4:O2	2.18	0.43
2:E:1174:LEU:HD23	2:E:1183:PHE:HD2	1.84	0.43
1:A:196:PHE:HA	1:A:206:VAL:O	2.19	0.42
1:B:196:PHE:CD2	1:B:205:VAL:CG1	2.96	0.42
1:B:204:VAL:HG22	2:E:1172:VAL:HB	2.01	0.42
1:B:750:VAL:HG21	1:B:911:MET:CE	2.49	0.42
1:A:297:GLU:HA	1:A:297:GLU:OE1	2.19	0.42
1:A:176:LEU:HD23	1:A:179:THR:OG1	2.19	0.42
1:A:831:GLY:HA2	4:A:990:SO4:O3	2.18	0.42
1:B:746:MET:O	1:B:897:VAL:CG2	2.67	0.42
1:A:807:MET:HE1	1:A:823:ILE:HD11	2.01	0.42
1:A:366:TYR:OH	1:A:370:LEU:HB2	2.19	0.42
1:B:499:THR:HG22	1:B:533:LYS:HG3	2.01	0.42
1:A:770:ALA:CB	1:A:803:LEU:HD13	2.50	0.42
1:B:753:GLY:HA3	1:B:768:THR:HB	2.00	0.42
1:B:330:THR:HG23	1:B:406:THR:HG22	2.02	0.41
1:B:242:LEU:HA	1:B:242:LEU:HD12	1.86	0.41
1:B:279:ASP:O	1:B:280:LEU:HD23	2.20	0.41
1:A:747:PHE:HB3	1:A:897:VAL:HG22	2.02	0.41
1:B:258:THR:HG23	1:B:261:GLN:HE21	1.86	0.41
1:A:169:THR:HG22	1:A:170:LEU:O	2.21	0.41
1:B:885:TRP:CZ3	1:B:888:LEU:HD23	2.56	0.41
1:B:876:PRO:HG2	1:B:911:MET:HE1	2.02	0.41
1:B:762:THR:HG23	1:B:762:THR:O	2.12	0.41
1:B:333:TRP:CD2	1:B:403:VAL:HG22	2.56	0.41
1:B:807:MET:CE	1:B:823:ILE:HD11	2.51	0.41
2:E:1190:MET:HA	2:E:1195:ALA:HB1	2.03	0.41
1:A:751:THR:CG2	1:A:768:THR:HG22	2.51	0.41
1:B:876:PRO:HG2	1:B:911:MET:CE	2.51	0.41
1:A:888:LEU:CD1	1:A:892:TYR:CE1	3.04	0.41
1:B:751:THR:HG23	1:B:768:THR:HG22	2.03	0.41
1:A:859:LEU:HD23	1:A:884:LEU:CD2	2.51	0.41
1:B:340:LYS:HD2	1:B:394:GLU:OE2	2.20	0.41
1:B:204:VAL:HG12	1:B:205:VAL:N	2.36	0.41
1:B:565:ILE:HD11	1:B:603:GLU:CG	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:ALA:HB3	1:A:431:ALA:HA	2.02	0.41
1:A:581:LYS:O	1:A:584:THR:O	2.39	0.40
1:B:403:VAL:CG1	1:B:403:VAL:O	2.68	0.40
2:D:1169:MET:CB	2:D:1190:MET:HE1	2.49	0.40
1:A:261:GLN:OE1	1:A:277:LEU:HD21	2.22	0.40
1:B:364:LEU:HD12	1:B:364:LEU:HA	1.85	0.40
1:A:580:LEU:O	1:A:584:THR:HB	2.21	0.40
1:A:298:ASP:OD2	1:A:300:TYR:HB2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	760/800 (95%)	728 (96%)	27 (4%)	5 (1%)	26	59
1	B	754/800 (94%)	720 (96%)	32 (4%)	2 (0%)	46	76
2	D	48/97 (50%)	42 (88%)	6 (12%)	0	100	100
2	E	56/97 (58%)	53 (95%)	3 (5%)	0	100	100
All	All	1618/1794 (90%)	1543 (95%)	68 (4%)	7 (0%)	39	71

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	759	SER
1	B	215	SER
1	A	216	GLN
1	A	339	LYS
1	A	287	GLU
1	A	586	GLU
1	B	175	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	673/699 (96%)	639 (95%)	34 (5%)	29	62
1	B	671/699 (96%)	629 (94%)	42 (6%)	22	50
2	D	49/82 (60%)	46 (94%)	3 (6%)	23	52
2	E	54/82 (66%)	46 (85%)	8 (15%)	4	10
All	All	1447/1562 (93%)	1360 (94%)	87 (6%)	24	53

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	128	ILE
1	A	156	ILE
1	A	179	THR
1	A	210	ARG
1	A	217	SER
1	A	267	GLU
1	A	271	GLU
1	A	276	THR
1	A	346	THR
1	A	356	LEU
1	A	370	LEU
1	A	404	GLU
1	A	405	VAL
1	A	425	SER
1	A	457	CYS
1	A	471	LEU
1	A	482	VAL
1	A	543	ARG
1	A	549	ARG
1	A	559	LEU
1	A	631	ARG
1	A	646	CYS
1	A	671	MET
1	A	684	LEU

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Mol	Chain	Res	Type
1	A	691	LEU
1	A	699	GLN
1	A	721	LEU
1	A	744	LYS
1	A	774	GLU
1	A	819	GLN
1	A	820	GLU
1	A	857	ARG
1	A	870	VAL
1	A	897	VAL
1	B	117	ASP
1	B	128	ILE
1	B	156	ILE
1	B	179	THR
1	B	205	VAL
1	B	209	CYS
1	B	210	ARG
1	B	215	SER
1	B	241	TRP
1	B	253	ARG
1	B	260	GLN
1	B	267	GLU
1	B	271	GLU
1	B	276	THR
1	B	279	ASP
1	B	357	MET
1	B	398	SER
1	B	404	GLU
1	B	405	VAL
1	B	425	SER
1	B	457	CYS
1	B	471	LEU
1	B	482	VAL
1	B	543	ARG
1	B	549	ARG
1	B	559	LEU
1	B	584	THR
1	B	631	ARG
1	B	684	LEU
1	B	691	LEU
1	B	699	GLN
1	B	721	LEU

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Mol	Chain	Res	Type
1	B	730	ARG
1	B	738	GLN
1	B	749	TYR
1	B	762	THR
1	B	774	GLU
1	B	819	GLN
1	B	820	GLU
1	B	857	ARG
1	B	870	VAL
1	B	897	VAL
2	D	1124	ASN
2	D	1125	LEU
2	D	1168	THR
2	E	1121	MET
2	E	1122	LEU
2	E	1125	LEU
2	E	1129	SER
2	E	1168	THR
2	E	1175	THR
2	E	1179	ASN
2	E	1197	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	ASN
1	A	146	ASN
1	A	150	ASN
1	A	185	ASN
1	A	304	ASN
1	A	328	ASN
1	A	408	ASN
1	A	443	HIS
1	A	458	GLN
1	A	524	ASN
1	A	629	GLN
1	A	665	GLN
1	A	699	GLN
1	A	722	GLN
1	A	741	GLN
1	A	860	ASN
1	A	886	ASN

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Mol	Chain	Res	Type
1	B	138	ASN
1	B	146	ASN
1	B	150	ASN
1	B	185	ASN
1	B	228	GLN
1	B	291	HIS
1	B	304	ASN
1	B	328	ASN
1	B	408	ASN
1	B	443	HIS
1	B	458	GLN
1	B	524	ASN
1	B	629	GLN
1	B	665	GLN
1	B	699	GLN
1	B	722	GLN
1	B	738	GLN
1	B	741	GLN
1	B	752	GLN
1	B	860	ASN
1	B	886	ASN

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](#)

Of 29 ligands modelled in this entry, 6 are monoatomic - leaving 23 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link

column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	A	987	-	4,4,4	0.12	0	6,6,6	0.10	0
4	SO4	A	988	-	4,4,4	0.23	0	6,6,6	0.22	0
4	SO4	A	989	-	4,4,4	0.24	0	6,6,6	0.32	0
4	SO4	A	990	-	4,4,4	0.39	0	6,6,6	0.71	0
4	SO4	A	991	-	4,4,4	0.18	0	6,6,6	0.13	0
4	SO4	A	992	-	4,4,4	0.31	0	6,6,6	0.37	0
4	SO4	A	993	-	4,4,4	0.14	0	6,6,6	0.26	0
4	SO4	A	994	-	4,4,4	0.29	0	6,6,6	0.34	0
4	SO4	A	995	-	4,4,4	0.26	0	6,6,6	0.12	0
4	SO4	A	996	-	4,4,4	0.23	0	6,6,6	0.19	0
4	SO4	A	997	-	4,4,4	0.20	0	6,6,6	0.16	0
4	SO4	A	998	-	4,4,4	0.24	0	6,6,6	0.18	0
4	SO4	A	999	-	4,4,4	0.25	0	6,6,6	0.14	0
4	SO4	B	989	-	4,4,4	0.20	0	6,6,6	0.40	0
4	SO4	B	990	-	4,4,4	0.09	0	6,6,6	0.43	0
4	SO4	B	991	-	4,4,4	0.22	0	6,6,6	0.18	0
4	SO4	B	992	-	4,4,4	0.12	0	6,6,6	0.13	0
4	SO4	B	993	-	4,4,4	0.15	0	6,6,6	0.09	0
4	SO4	B	995	-	4,4,4	0.22	0	6,6,6	0.26	0
4	SO4	B	996	-	4,4,4	0.29	0	6,6,6	0.13	0
4	SO4	B	997	-	4,4,4	0.23	0	6,6,6	0.16	0
4	SO4	B	998	-	4,4,4	0.13	0	6,6,6	0.22	0
4	SO4	B	999	-	4,4,4	0.37	0	6,6,6	0.61	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
4	SO4	A	987	-	-	0/0/0/0	0/0/0/0
4	SO4	A	988	-	-	0/0/0/0	0/0/0/0
4	SO4	A	989	-	-	0/0/0/0	0/0/0/0
4	SO4	A	990	-	-	0/0/0/0	0/0/0/0
4	SO4	A	991	-	-	0/0/0/0	0/0/0/0
4	SO4	A	992	-	-	0/0/0/0	0/0/0/0
4	SO4	A	993	-	-	0/0/0/0	0/0/0/0
4	SO4	A	994	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SO4	A	995	-	-	0/0/0/0	0/0/0/0
4	SO4	A	996	-	-	0/0/0/0	0/0/0/0
4	SO4	A	997	-	-	0/0/0/0	0/0/0/0
4	SO4	A	998	-	-	0/0/0/0	0/0/0/0
4	SO4	A	999	-	-	0/0/0/0	0/0/0/0
4	SO4	B	989	-	-	0/0/0/0	0/0/0/0
4	SO4	B	990	-	-	0/0/0/0	0/0/0/0
4	SO4	B	991	-	-	0/0/0/0	0/0/0/0
4	SO4	B	992	-	-	0/0/0/0	0/0/0/0
4	SO4	B	993	-	-	0/0/0/0	0/0/0/0
4	SO4	B	995	-	-	0/0/0/0	0/0/0/0
4	SO4	B	996	-	-	0/0/0/0	0/0/0/0
4	SO4	B	997	-	-	0/0/0/0	0/0/0/0
4	SO4	B	998	-	-	0/0/0/0	0/0/0/0
4	SO4	B	999	-	-	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.

9 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	988	SO4	1	0
4	A	990	SO4	1	0
4	A	994	SO4	2	0
4	A	996	SO4	1	0
4	A	998	SO4	2	0
4	A	999	SO4	1	0
4	B	991	SO4	1	0
4	B	996	SO4	1	0
4	B	999	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 ⓘ

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	772/800 (96%)	0.22	35 (4%) 37 31	9, 20, 33, 44	0
1	B	768/800 (96%)	0.16	33 (4%) 39 32	8, 20, 34, 46	0
2	D	54/97 (55%)	0.35	2 (3%) 45 38	9, 29, 45, 47	0
2	E	60/97 (61%)	0.64	10 (16%) 2 1	9, 34, 49, 52	0
All	All	1654/1794 (92%)	0.21	80 (4%) 34 28	8, 20, 38, 52	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	1179	ASN	6.3
1	A	377	ILE	5.3
1	A	337	LEU	5.1
1	A	368	GLY	4.7
1	A	373	LEU	4.6
2	E	1104	GLY	4.4
1	A	587	LEU	4.0
1	B	363	CYS	4.0
1	B	736	ASP	4.0
1	A	401	ALA	3.9
2	E	1181	GLN	3.9
1	B	399	VAL	3.6
2	E	1183	PHE	3.6
1	B	368	GLY	3.5
1	B	401	ALA	3.5
1	A	278	GLU	3.5
1	B	404	GLU	3.4
1	B	733	LYS	3.4
1	A	734	GLY	3.4
1	A	338	ASN	3.3
1	B	456	LYS	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	217	SER	3.2
1	A	339	LYS	3.2
1	A	372	PRO	3.1
1	A	736	ASP	3.1
2	E	1178	GLY	3.1
2	E	1182	GLN	3.0
1	B	398	SER	3.0
1	B	734	GLY	2.9
1	B	120	ILE	2.9
1	B	846	GLU	2.9
1	B	376	GLY	2.9
1	A	361	GLU	2.8
1	B	375	LYS	2.8
1	A	369	ASP	2.8
1	A	374	TRP	2.8
1	A	846	GLU	2.7
1	B	395	LEU	2.7
1	B	271	GLU	2.6
1	A	331	VAL	2.6
1	A	376	GLY	2.6
1	B	412	ASP	2.6
1	A	405	VAL	2.6
1	A	271	GLU	2.6
1	A	578	GLN	2.6
1	B	365	ARG	2.6
1	B	411	VAL	2.6
1	B	848	GLN	2.6
1	A	224	TRP	2.6
1	A	591	ASP	2.6
1	B	374	TRP	2.6
1	B	377	ILE	2.5
1	B	364	LEU	2.5
1	A	166	LYS	2.5
1	A	375	LYS	2.5
1	B	117	ASP	2.4
1	A	366	TYR	2.4
2	E	1198	HIS	2.4
1	A	367	LYS	2.4
1	B	397	SER	2.4
1	B	341	ARG	2.4
2	D	1183	PHE	2.4
1	A	362	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
2	D	1127	GLN	2.2
1	B	413	PHE	2.2
1	B	175	PRO	2.2
1	A	173	ASP	2.2
1	A	356	LEU	2.1
1	A	363	CYS	2.1
1	A	342	ILE	2.1
1	B	743	ASP	2.1
1	B	849	GLY	2.1
1	B	342	ILE	2.1
2	E	1176	ARG	2.1
1	B	592	GLU	2.1
2	E	1184	LYS	2.1
2	E	1103	MET	2.0
1	A	588	SER	2.0
1	B	356	LEU	2.0
1	A	371	ALA	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
4	SO4	A	997	5/5	0.91	0.35	6.72	133,133,133,133	0
4	SO4	A	999	5/5	0.93	0.41	4.26	135,135,135,136	0
4	SO4	B	999	5/5	0.88	0.22	2.44	73,75,76,76	0
4	SO4	B	996	5/5	0.89	0.32	2.37	152,152,153,153	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	SO4	A	996	5/5	0.81	0.32	2.01	138,138,138,138	0
4	SO4	B	997	5/5	0.89	0.25	0.59	131,131,131,131	0
4	SO4	A	992	5/5	0.83	0.21	-0.37	96,97,97,97	0
4	SO4	A	990	5/5	0.94	0.17	-0.52	71,72,73,73	0
3	ZN	B	3	1/1	0.99	0.13	-0.79	13,13,13,13	0
3	ZN	B	1	1/1	0.99	0.12	-1.00	10,10,10,10	0
4	SO4	B	991	5/5	0.88	0.20	-1.21	151,151,152,152	0
3	ZN	A	1	1/1	1.00	0.10	-1.25	11,11,11,11	0
4	SO4	A	991	5/5	0.92	0.10	-1.30	117,118,118,118	0
4	SO4	A	995	5/5	0.95	0.13	-1.41	96,96,97,97	0
3	ZN	A	3	1/1	0.98	0.10	-1.46	19,19,19,19	0
4	SO4	B	995	5/5	0.95	0.13	-1.83	79,80,81,82	0
4	SO4	B	998	5/5	0.87	0.18	-1.98	104,104,105,105	0
3	ZN	A	2	1/1	0.99	0.06	-2.14	20,20,20,20	1
3	ZN	B	2	1/1	0.99	0.08	-2.80	24,24,24,24	1
4	SO4	A	989	5/5	0.89	0.17	-	113,114,114,114	0
4	SO4	A	988	5/5	0.79	0.22	-	141,141,142,142	0
4	SO4	A	994	5/5	0.84	0.16	-	126,126,126,127	0
4	SO4	B	993	5/5	0.93	0.12	-	106,107,107,108	0
4	SO4	B	992	5/5	0.76	0.25	-	129,129,129,129	0
4	SO4	A	987	5/5	0.73	0.27	-	127,127,127,128	0
4	SO4	B	990	5/5	0.81	0.31	-	110,111,111,112	0
4	SO4	A	993	5/5	0.88	0.14	-	113,113,113,113	0
4	SO4	A	998	5/5	0.87	0.18	-	119,120,120,120	0
4	SO4	B	989	5/5	0.76	0.24	-	118,118,119,119	0

6.5 Other polymers ⓘ

There are no such residues in this entry.