



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:53 PM UTC

PDB ID : 4YJK / pdb_00004yjk
Title : Crystal structure of C212S mutant of Shewanella oneidensis MR-1 uridine phosphorylase
Authors : Safonova, T.N.; Mordkovich, N.N.; Manuvera, V.A.; Dorovatovsky, P.V.; Veiko, V.P.; Popov, V.O.; Polyakov, K.M.
Deposited on : 2015-03-03
Resolution : 1.68 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

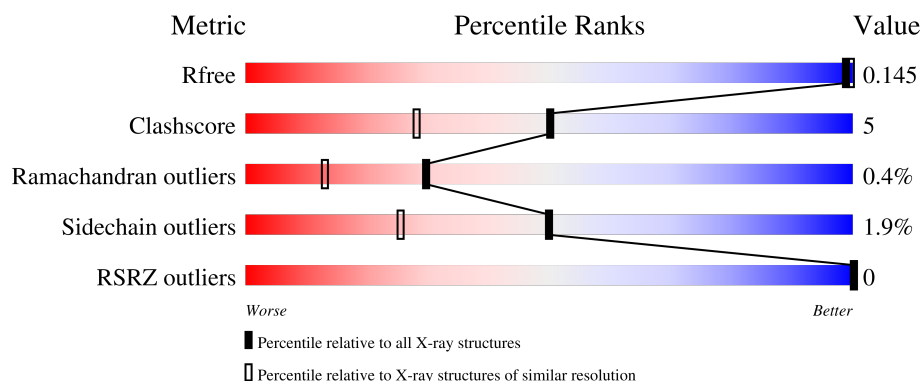
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

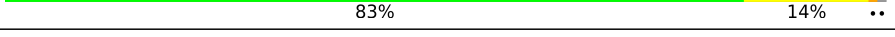
The reported resolution of this entry is 1.68 Å.

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}	180053	1054 (1.68-1.68)
Clashscore	190562	1078 (1.68-1.68)
Ramachandran outliers	187476	1068 (1.68-1.68)
Sidechain outliers	187428	1067 (1.68-1.68)
RSRZ outliers	180081	1055 (1.68-1.68)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	252	
1	B	252	
1	C	252	
1	D	252	
1	E	252	

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Mol	Chain	Length	Quality of chain
1	F	252	 84% 11% . .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	D	303	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11842 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.

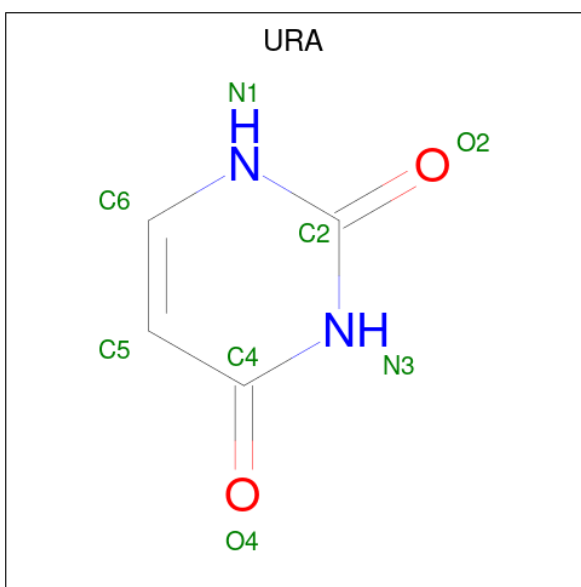
- Molecule 1 is a protein called Uridine phosphorylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	249	Total	C	N	O	S	0	2	0
			1851	1160	317	361	13			
1	F	244	Total	C	N	O	S	0	1	0
			1817	1141	312	352	12			
1	A	245	Total	C	N	O	S	0	0	0
			1816	1140	313	351	12			
1	C	249	Total	C	N	O	S	0	2	0
			1859	1166	320	360	13			
1	E	242	Total	C	N	O	S	0	1	0
			1796	1128	311	345	12			
1	B	240	Total	C	N	O	S	0	1	0
			1783	1119	310	342	12			

There are 6 discrepancies between the modelled and reference sequences:

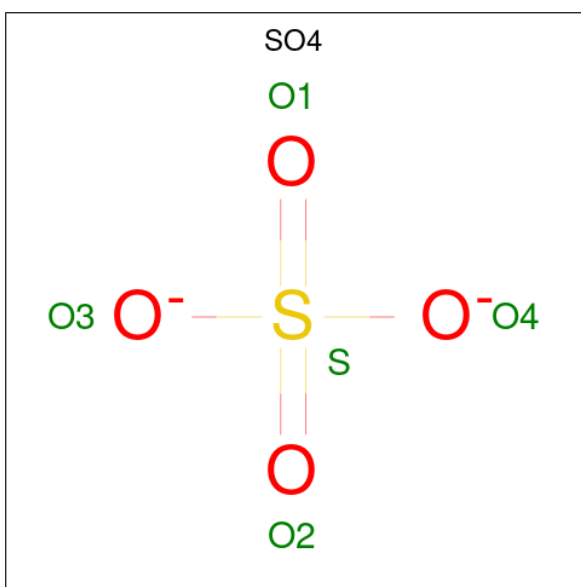
Chain	Residue	Modelled	Actual	Comment	Reference
D	212	SER	CYS	engineered mutation	UNP Q8E9X9
F	212	SER	CYS	engineered mutation	UNP Q8E9X9
A	212	SER	CYS	engineered mutation	UNP Q8E9X9
C	212	SER	CYS	engineered mutation	UNP Q8E9X9
E	212	SER	CYS	engineered mutation	UNP Q8E9X9
B	212	SER	CYS	engineered mutation	UNP Q8E9X9

- Molecule 2 is URACIL (CCD ID: URA) (formula: $C_4H_4N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	N	O	0	0
			8	4	2	2		
2	F	1	Total	C	N	O	0	0
			8	4	2	2		
2	A	1	Total	C	N	O	0	0
			8	4	2	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	1
			10	8	2		
3	F	1	Total	O	S	0	1
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	1
			10	8	2		
3	E	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

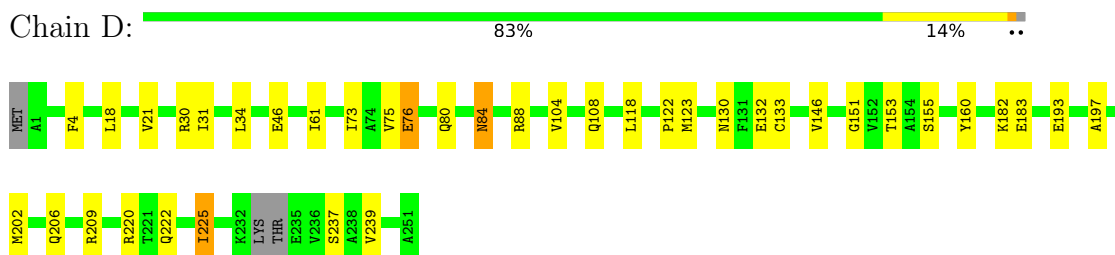
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	149	Total	O	0	0
			149	149		
4	F	138	Total	O	0	0
			138	138		
4	A	151	Total	O	0	0
			151	151		
4	C	149	Total	O	0	0
			149	149		
4	E	129	Total	O	0	0
			129	129		
4	B	125	Total	O	0	0
			125	125		

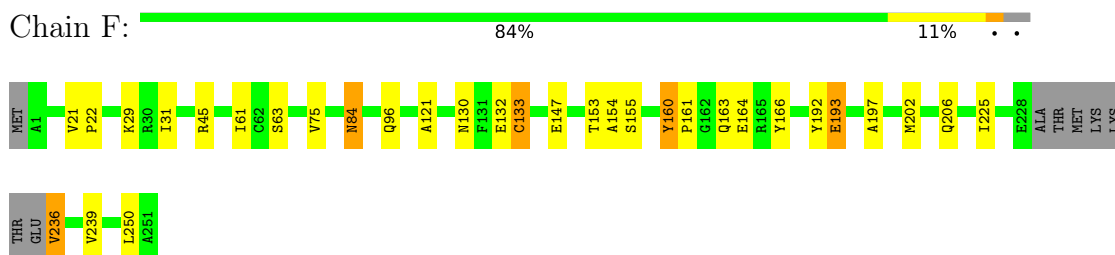
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

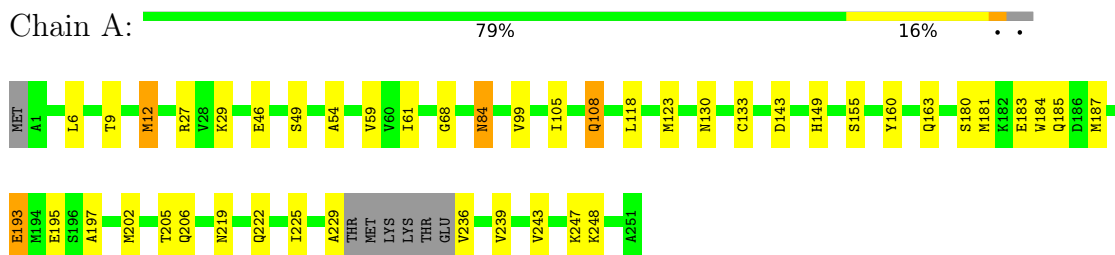
- Molecule 1: Uridine phosphorylase



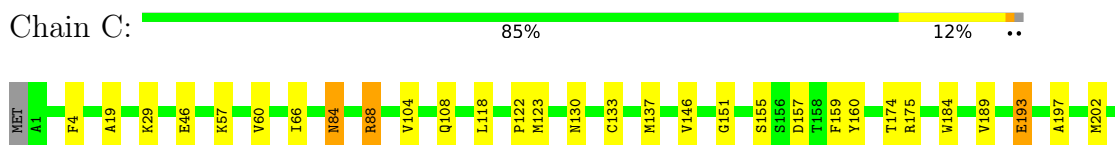
- Molecule 1: Uridine phosphorylase

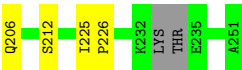


- Molecule 1: Uridine phosphorylase

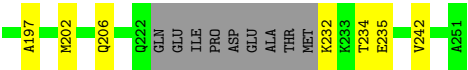
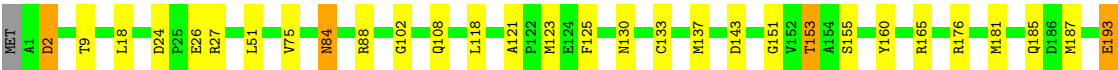
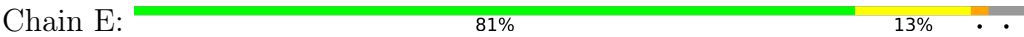


- Molecule 1: Uridine phosphorylase

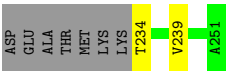
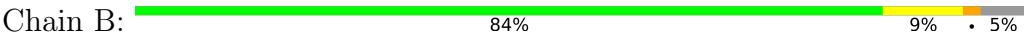




● Molecule 1: Uridine phosphorylase



● Molecule 1: Uridine phosphorylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.23Å 95.40Å 91.20Å 90.00° 119.98° 90.00°	Depositor
Resolution (Å)	19.35 – 1.68 19.35 – 1.68	Depositor EDS
% Data completeness (in resolution range)	98.2 (19.35-1.68) 98.3 (19.35-1.68)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 1.68Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.138 , 0.161 (Not available) , 0.145	Depositor DCC
R_{free} test set	7594 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	10.8	Xtriage
Anisotropy	0.818	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 20.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.35$, $\langle L^2 \rangle = 0.18$	Xtriage
Estimated twinning fraction	0.327 for l,k,-h-l 0.327 for -h-l,k,h 0.257 for -h-l,-k,l 0.286 for h,-k,-h-l 0.246 for l,-k,h	Xtriage
Reported twinning fraction	0.471 for H, K, L 0.064 for H+L, -K, -L 0.056 for L, -K, H 0.060 for L, K, -H-L 0.182 for -H-L, K, H 0.166 for -H, -K, H+L	Depositor
Outliers	0 of 151174 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11842	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.72% 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, 0.333 respectively for untwinned datasets, and 0.375, 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, URA

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	1.23	2/1848 (0.1%)	1.27	6/2514 (0.2%)
1	B	1.20	6/1818 (0.3%)	1.17	3/2471 (0.1%)
1	C	1.24	4/1899 (0.2%)	1.25	5/2582 (0.2%)
1	D	1.24	4/1892 (0.2%)	1.26	4/2573 (0.2%)
1	E	1.24	3/1831 (0.2%)	1.19	5/2490 (0.2%)
1	F	1.22	0/1854	1.22	11/2522 (0.4%)
All	All	1.23	19/11142 (0.2%)	1.23	34/15152 (0.2%)

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	184	TRP	C-O	-6.67	1.16	1.24
1	A	59	VAL	C-O	-6.08	1.18	1.24
1	E	75	VAL	C-O	-5.88	1.16	1.24
1	E	51	LEU	N-CA	-5.62	1.39	1.46
1	B	111	VAL	C-O	5.61	1.30	1.24
1	D	76	GLU	C-O	5.61	1.30	1.24
1	B	201	THR	N-CA	-5.46	1.39	1.46
1	C	157	ASP	CA-C	5.43	1.60	1.52
1	C	184	TRP	N-CA	5.41	1.52	1.46
1	B	155	SER	C-O	5.38	1.30	1.23
1	C	137	MET	C-O	-5.32	1.17	1.24
1	E	153	THR	N-CA	5.32	1.52	1.46
1	B	68	GLY	C-O	-5.20	1.18	1.24
1	B	108	GLN	CA-C	5.19	1.59	1.52
1	D	122	PRO	N-CA	-5.11	1.40	1.47
1	D	76	GLU	N-CA	5.11	1.52	1.46
1	D	220	ARG	C-O	-5.06	1.17	1.24
1	B	78	LEU	C-O	-5.03	1.18	1.24
1	C	189	VAL	C-O	5.01	1.29	1.24

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	68	GLY	N-CA-C	7.34	124.55	115.22
1	D	75	VAL	N-CA-C	-7.34	103.63	110.53
1	E	75	VAL	N-CA-C	-6.90	103.48	111.00
1	B	119	HIS	N-CA-C	-6.53	105.06	113.16
1	F	225	ILE	CB-CA-C	6.50	117.42	110.53
1	A	163	GLN	N-CA-C	-6.36	104.67	112.88
1	F	75	VAL	N-CA-C	-6.32	104.59	110.53
1	A	243	VAL	N-CA-C	-6.31	104.12	111.00
1	F	133	CYS	N-CA-C	-6.14	104.49	112.23
1	D	225	ILE	CA-C-N	-5.99	113.80	120.14
1	D	225	ILE	C-N-CA	-5.99	113.80	120.14
1	E	121	ALA	O-C-N	5.92	126.45	121.83
1	D	222	GLN	N-CA-C	5.85	121.30	114.04
1	E	2	ASP	N-CA-C	-5.68	104.70	111.69
1	C	57	LYS	CA-C-N	-5.68	114.08	119.76
1	C	57	LYS	C-N-CA	-5.68	114.08	119.76
1	C	4	PHE	N-CA-C	5.67	118.40	111.82
1	F	121	ALA	O-C-N	5.64	126.23	121.83
1	A	6	LEU	N-CA-C	5.62	118.34	111.82
1	B	163	GLN	N-CA-C	-5.57	105.70	112.88
1	F	163	GLN	N-CA-C	-5.49	105.89	112.59
1	A	12	MET	N-CA-C	5.38	118.64	111.75
1	F	21	VAL	CA-C-N	-5.26	115.32	120.52
1	F	21	VAL	C-N-CA	-5.26	115.32	120.52
1	B	197	ALA	N-CA-C	-5.23	105.47	111.07
1	F	96	GLN	CA-C-N	-5.12	114.39	119.56
1	F	96	GLN	C-N-CA	-5.12	114.39	119.56
1	A	108	GLN	CB-CA-C	-5.11	102.17	110.85
1	E	125	PHE	CA-C-N	-5.11	114.73	120.14
1	E	125	PHE	C-N-CA	-5.11	114.73	120.14
1	C	66	ILE	N-CA-C	5.08	115.64	109.30
1	F	147	GLU	CA-C-N	-5.07	115.35	120.31
1	F	147	GLU	C-N-CA	-5.07	115.35	120.31
1	C	122	PRO	N-CA-C	-5.00	103.91	111.41

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	1816	0	1795	24	0
1	B	1783	0	1769	17	0
1	C	1859	0	1837	18	0
1	D	1851	0	1822	26	0
1	E	1796	0	1779	20	0
1	F	1817	0	1799	16	0
2	A	8	0	3	0	0
2	D	8	0	3	0	0
2	F	8	0	3	0	0
3	A	5	0	0	1	0
3	B	5	0	0	1	0
3	C	5	0	0	0	0
3	D	20	0	0	3	0
3	E	15	0	0	1	0
3	F	5	0	0	0	0
4	A	151	0	0	4	0
4	B	125	0	0	1	0
4	C	149	0	0	1	0
4	D	149	0	0	1	0
4	E	129	0	0	4	0
4	F	138	0	0	1	0
All	All	11842	0	10810	115	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 5.

All (115) 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:D:30:ARG:NH2	3:D:303:SO4:O1	2.05	0.89
1:D:237:SER:HB2	3:D:303:SO4:S	2.21	0.80
1:F:45:ARG:NH2	3:E:302:SO4:O2	2.16	0.79
1:E:27:ARG:NH2	4:E:515:HOH:O	2.18	0.76
1:E:143:ASP:OD1	4:E:401:HOH:O	2.06	0.71
1:A:108:GLN:NE2	4:A:468:HOH:O	2.24	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:130:ASN:HD22	1:E:133:CYS:H	1.38	0.69
1:A:155:SER:HB3	1:A:197:ALA:HB2	1.75	0.69
1:D:130:ASN:HD22	1:D:133:CYS:H	1.41	0.69
1:D:108:GLN:HE21	1:D:151:GLY:HA2	1.59	0.67
1:A:193:GLU:HB2	4:A:515:HOH:O	1.95	0.65
1:D:202:MET:O	1:D:206:GLN:HG2	1.99	0.63
1:A:143:ASP:OD2	1:A:248:LYS:NZ	2.28	0.63
1:D:84:ASN:C	1:D:84:ASN:HD22	2.08	0.61
1:A:229:ALA:HB3	4:A:525:HOH:O	2.00	0.61
1:A:130:ASN:HD22	1:A:133:CYS:H	1.49	0.61
1:C:84:ASN:HD22	1:C:84:ASN:C	2.08	0.60
1:C:130:ASN:HD22	1:C:133:CYS:H	1.48	0.60
1:B:130:ASN:HD22	1:B:133:CYS:H	1.48	0.60
1:D:155:SER:HB3	1:D:197:ALA:HB2	1.84	0.60
1:B:84:ASN:C	1:B:84:ASN:HD22	2.10	0.59
1:A:202:MET:O	1:A:206:GLN:HG2	2.03	0.59
1:E:165:ARG:HD2	4:E:511:HOH:O	2.04	0.58
1:F:130:ASN:HD22	1:F:133:CYS:H	1.52	0.58
1:C:202:MET:O	1:C:206:GLN:HG2	2.03	0.58
1:E:84:ASN:C	1:E:84:ASN:HD22	2.12	0.57
1:D:153:THR:HG21	1:D:193[A]:GLU:HG3	1.85	0.57
1:D:108:GLN:NE2	4:D:504:HOH:O	2.13	0.56
1:D:132:GLU:OE1	1:D:209:ARG:NH1	2.38	0.56
1:A:222:GLN:NE2	4:A:541:HOH:O	2.37	0.56
1:A:84:ASN:C	1:A:84:ASN:HD22	2.13	0.56
1:D:118:LEU:HD21	1:D:123:MET:HE2	1.88	0.56
1:E:108:GLN:HE21	1:E:151:GLY:HA2	1.71	0.55
1:E:176:ARG:HD2	4:E:439:HOH:O	2.05	0.55
1:D:153:THR:HG21	1:D:193[B]:GLU:HG3	1.88	0.55
1:B:102:GLY:HA2	1:B:234:THR:CG2	2.37	0.55
1:F:155:SER:HB3	1:F:197:ALA:HB2	1.91	0.53
1:B:27:ARG:NH2	3:B:301:SO4:O4	2.41	0.53
1:C:193[A]:GLU:HB2	4:C:512:HOH:O	2.08	0.53
1:C:19:ALA:HB2	1:C:60:VAL:HG13	1.90	0.53
1:C:118:LEU:HD21	1:C:123:MET:HE2	1.89	0.53
1:C:108:GLN:HE21	1:C:151:GLY:HA2	1.74	0.53
1:C:225:ILE:HG22	1:C:226:PRO:O	2.08	0.53
1:E:155:SER:HB3	1:E:197:ALA:HB2	1.90	0.53
1:F:202:MET:O	1:F:206:GLN:HG2	2.08	0.52
1:E:232:LYS:HA	1:E:235:GLU:HG3	1.91	0.52
1:B:202:MET:O	1:B:206:GLN:HG2	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:237:SER:HB2	3:D:303:SO4:O1	2.10	0.52
1:C:88[A]:ARG:HG2	1:C:212:SER:OG	2.10	0.52
1:C:155:SER:HB3	1:C:197:ALA:HB2	1.92	0.51
1:A:236:VAL:N	1:A:239:VAL:HG22	2.25	0.51
1:D:31:ILE:HG12	1:D:239:VAL:HG13	1.92	0.51
1:F:61:ILE:HD12	1:F:61:ILE:N	2.26	0.50
1:B:155:SER:HB3	1:B:197:ALA:HB2	1.93	0.50
1:D:104:VAL:HG22	1:D:146:VAL:HG11	1.94	0.50
1:C:19:ALA:HA	1:C:60:VAL:O	2.12	0.49
1:F:84:ASN:C	1:F:84:ASN:HD22	2.20	0.49
1:B:30:ARG:CD	4:B:522:HOH:O	2.61	0.49
1:E:137:MET:HE1	1:E:242:VAL:HG23	1.95	0.48
1:A:9:THR:O	1:A:12:MET:HB2	2.14	0.48
1:A:27:ARG:NH2	3:A:302:SO4:O2	2.42	0.48
1:F:22:PRO:O	1:F:63:SER:HA	2.13	0.48
1:E:130:ASN:ND2	1:E:133:CYS:H	2.11	0.47
1:F:31:ILE:HG12	1:F:239:VAL:HG13	1.97	0.47
1:B:102:GLY:HA2	1:B:234:THR:HG22	1.98	0.46
1:A:181:MET:O	1:A:185:GLN:HG3	2.16	0.46
1:D:4:PHE:CD1	1:C:225:ILE:HG23	2.50	0.45
1:D:46:GLU:HB3	1:C:46:GLU:HB3	1.98	0.45
1:A:54:ALA:HB1	1:A:247:LYS:HG2	1.98	0.45
1:A:183:GLU:O	1:A:187:MET:HG3	2.17	0.45
1:B:24:ASP:CB	1:B:27:ARG:HH21	2.29	0.45
1:A:99:VAL:O	1:A:219:ASN:ND2	2.50	0.44
1:A:118:LEU:HD21	1:A:123:MET:HE2	1.98	0.44
1:E:2:ASP:CG	1:E:9:THR:HA	2.43	0.44
1:B:34:LEU:HD12	1:B:239:VAL:HG12	2.00	0.44
1:E:24:ASP:OD1	1:E:26:GLU:HG3	2.18	0.44
1:E:153:THR:HG21	1:E:193:GLU:HG3	2.00	0.43
1:D:73:ILE:HD11	1:C:159:PHE:HB2	1.99	0.43
1:D:34:LEU:CD1	1:D:239:VAL:HG12	2.49	0.43
1:A:49:SER:HA	1:A:61:ILE:O	2.18	0.43
1:C:104:VAL:HG22	1:C:146:VAL:HG11	2.01	0.43
1:C:174:THR:O	1:C:175:ARG:C	2.62	0.42
1:D:130:ASN:ND2	1:D:133:CYS:H	2.11	0.42
1:E:187:MET:HE1	1:B:122:PRO:HD2	2.00	0.42
1:B:19:ALA:HB2	1:B:60:VAL:HG13	2.01	0.42
1:E:18:LEU:C	1:E:18:LEU:HD23	2.44	0.42
1:A:130:ASN:ND2	1:A:133:CYS:H	2.15	0.42
1:D:21:VAL:HG23	1:D:88:ARG:HA	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:SER:O	1:A:183:GLU:HB2	2.20	0.42
1:E:102:GLY:HA2	1:E:234:THR:OG1	2.19	0.42
1:E:202:MET:O	1:E:206:GLN:HG2	2.19	0.42
1:D:18:LEU:C	1:D:18:LEU:HD23	2.45	0.42
1:E:118:LEU:HD21	1:E:123:MET:HE2	2.00	0.42
1:D:153:THR:CG2	1:D:193[A]:GLU:HG3	2.50	0.42
1:F:164:GLU:HG2	1:F:166:TYR:CE2	2.55	0.41
1:C:19:ALA:CB	1:C:60:VAL:HG13	2.50	0.41
1:B:18:LEU:C	1:B:18:LEU:HD23	2.45	0.41
1:D:182:LYS:O	1:D:183:GLU:C	2.63	0.41
1:B:102:GLY:HA2	1:B:234:THR:HG21	2.01	0.41
1:F:154:ALA:HB3	1:F:192:TYR:CD1	2.55	0.41
1:A:46:GLU:HB3	1:B:46:GLU:HB3	2.01	0.41
1:F:236:VAL:HA	4:F:524:HOH:O	2.21	0.41
1:B:130:ASN:ND2	1:B:133:CYS:H	2.14	0.41
1:A:105:ILE:HG12	1:A:149:HIS:HB2	2.01	0.41
1:C:130:ASN:ND2	1:C:133:CYS:H	2.17	0.41
1:F:160:TYR:CB	1:F:161:PRO:CD	2.99	0.41
1:F:160:TYR:HB2	1:F:161:PRO:CD	2.50	0.41
1:A:193:GLU:CD	1:A:195:GLU:H	2.29	0.41
1:F:130:ASN:HD21	1:F:132:GLU:HB3	1.86	0.41
1:F:153:THR:HG21	1:F:193[A]:GLU:HG3	2.03	0.41
1:E:181:MET:O	1:E:185:GLN:HG3	2.21	0.41
1:D:76:GLU:O	1:D:80:GLN:HG3	2.20	0.40
1:F:29:LYS:HB3	1:F:29:LYS:HE3	1.88	0.40
1:A:205:THR:O	1:B:172:ARG:NH2	2.55	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	241/252 (96%)	234 (97%)	6 (2%)	1 (0%)	30	13
1	B	237/252 (94%)	233 (98%)	3 (1%)	1 (0%)	30	13
1	C	247/252 (98%)	240 (97%)	6 (2%)	1 (0%)	30	13
1	D	247/252 (98%)	243 (98%)	3 (1%)	1 (0%)	30	13
1	E	239/252 (95%)	234 (98%)	4 (2%)	1 (0%)	30	13
1	F	241/252 (96%)	234 (97%)	6 (2%)	1 (0%)	30	13
All	All	1452/1512 (96%)	1418 (98%)	28 (2%)	6 (0%)	30	13

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	160	TYR
1	A	160	TYR
1	C	160	TYR
1	B	160	TYR
1	F	160	TYR
1	E	160	TYR

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	189/198 (96%)	185 (98%)	4 (2%)	47	21
1	B	186/198 (94%)	181 (97%)	5 (3%)	39	13
1	C	195/198 (98%)	189 (97%)	6 (3%)	35	10
1	D	194/198 (98%)	192 (99%)	2 (1%)	68	52
1	E	187/198 (94%)	183 (98%)	4 (2%)	47	21
1	F	191/198 (96%)	186 (97%)	5 (3%)	40	15
All	All	1142/1188 (96%)	1116 (98%)	26 (2%)	50	18

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

Mol	Chain	Res	Type
1	D	84	ASN
1	D	225	ILE
1	F	84	ASN
1	F	193[A]	GLU
1	F	193[B]	GLU
1	F	236	VAL
1	F	250	LEU
1	A	29	LYS
1	A	84	ASN
1	A	193	GLU
1	A	225	ILE
1	C	29	LYS
1	C	84	ASN
1	C	88[A]	ARG
1	C	88[B]	ARG
1	C	193[A]	GLU
1	C	193[B]	GLU
1	E	84	ASN
1	E	88[A]	ARG
1	E	88[B]	ARG
1	E	193	GLU
1	B	27	ARG
1	B	30	ARG
1	B	84	ASN
1	B	108	GLN
1	B	193	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	84	ASN
1	D	98	HIS
1	D	108	GLN
1	D	130	ASN
1	F	84	ASN
1	F	130	ASN
1	F	222	GLN
1	F	223	GLN
1	A	37	ASN
1	A	84	ASN
1	A	98	HIS
1	A	108	GLN
1	A	130	ASN

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Mol	Chain	Res	Type
1	A	223	GLN
1	C	37	ASN
1	C	84	ASN
1	C	108	GLN
1	C	130	ASN
1	E	84	ASN
1	E	108	GLN
1	E	130	ASN
1	B	84	ASN
1	B	108	GLN
1	B	130	ASN
1	B	191	ASN
1	B	206	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

14 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	D	304[A]	-	4,4,4	0.48	0	6,6,6	0.21	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	E	301[B]	-	4,4,4	0.44	0	6,6,6	0.35	0
3	SO4	E	302	-	4,4,4	0.44	0	6,6,6	0.74	0
2	URA	A	301	-	8,8,8	0.68	0	10,10,10	0.99	1 (10%)
3	SO4	B	301	-	4,4,4	0.60	0	6,6,6	0.33	0
2	URA	D	301	-	8,8,8	0.89	0	10,10,10	1.08	1 (10%)
3	SO4	A	302	-	4,4,4	0.75	0	6,6,6	0.45	0
2	URA	F	301	-	8,8,8	0.85	0	10,10,10	1.05	1 (10%)
3	SO4	D	302	-	4,4,4	0.81	0	6,6,6	0.38	0
3	SO4	D	303	-	4,4,4	0.37	0	6,6,6	0.78	0
3	SO4	C	301	-	4,4,4	0.46	0	6,6,6	0.60	0
3	SO4	D	304[B]	-	4,4,4	0.38	0	6,6,6	0.73	0
3	SO4	E	301[A]	-	4,4,4	0.41	0	6,6,6	0.39	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	URA	F	301	-	-	-	0/1/1/1
2	URA	A	301	-	-	-	0/1/1/1
2	URA	D	301	-	-	-	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	301	URA	C6-N1-C2	-2.54	120.84	122.40
2	D	301	URA	C6-N1-C2	-2.45	120.90	122.40
2	A	301	URA	C6-N1-C2	-2.19	121.06	122.40

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	302	SO4	1	0
3	B	301	SO4	1	0
3	A	302	SO4	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	303	SO4	3	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	245/252 (97%)	-1.23	0 100 100	8, 13, 24, 39	0
1	B	240/252 (95%)	-1.19	0 100 100	8, 14, 28, 37	1 (0%)
1	C	249/252 (98%)	-1.23	0 100 100	7, 13, 23, 41	2 (0%)
1	D	249/252 (98%)	-1.23	0 100 100	6, 13, 23, 42	2 (0%)
1	E	242/252 (96%)	-1.23	0 100 100	8, 13, 24, 41	1 (0%)
1	F	244/252 (96%)	-1.24	0 100 100	7, 13, 24, 31	1 (0%)
All	All	1469/1512 (97%)	-1.22	0 100 100	6, 13, 25, 42	7 (0%)

There are no RSRZ outliers to report.

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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
3	SO4	D	303	5/5	0.98	0.06	23,23,24,25	5

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	E	302	5/5	0.98	0.07	21,21,23,23	5
3	SO4	B	301	5/5	0.98	0.05	18,20,21,21	5
2	URA	D	301	8/8	0.99	0.04	13,14,14,15	0
3	SO4	D	304[A]	5/5	0.99	0.03	17,18,18,18	5
3	SO4	D	304[B]	5/5	0.99	0.03	14,14,15,15	5
3	SO4	A	302	5/5	0.99	0.05	16,17,19,20	0
3	SO4	C	301	5/5	0.99	0.04	13,14,14,15	5
3	SO4	E	301[A]	5/5	0.99	0.03	17,18,19,20	5
3	SO4	E	301[B]	5/5	0.99	0.03	16,16,17,18	5
2	URA	F	301	8/8	0.99	0.03	12,13,13,15	0
2	URA	A	301	8/8	0.99	0.03	14,16,17,19	0
3	SO4	D	302	5/5	1.00	0.03	17,17,18,19	0
3	SO4	F	302[A]	5/5	1.00	0.03	13,14,14,16	1

6.5 Other polymers [i](#)

There are no such residues in this entry.