



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 6, 2024 – 06:11 PM JST

PDB ID : 5XVC
Title : [NiFe]-hydrogenase (Hyb-type) from *Citrobacter* sp. S-77 in a ferricyanide-oxidized condition
Authors : Nishikawa, K.; Matsuura, H.; Muhd Noor, N.D.; Tai, H.; Hirota, S.; Kim, J.; Kang, J.; Tateno, M.; Yoon, K.S.; Ogo, S.; Shomura, Y.; Higuchi, Y.
Deposited on : 2017-06-27
Resolution : 2.05 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	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

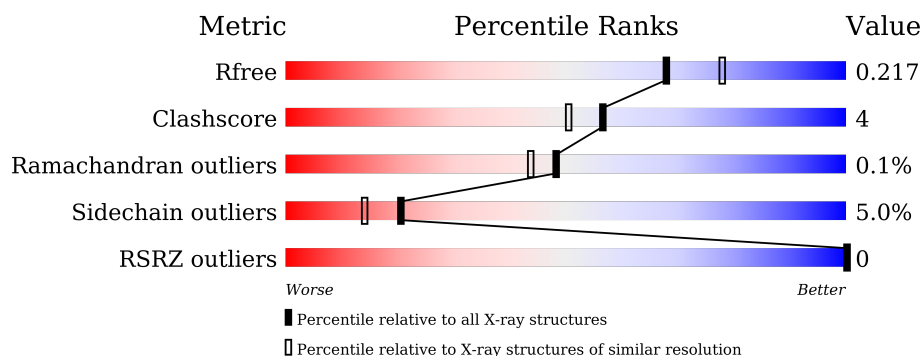
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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}	164625	2096 (2.04-2.04)
Clashscore	180529	2229 (2.04-2.04)
Ramachandran outliers	177936	2217 (2.04-2.04)
Sidechain outliers	177891	2217 (2.04-2.04)
RSRZ outliers	164620	2096 (2.04-2.04)

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	L	552	<div> <div style="width: 89%;"></div> <div style="width: 10%;"></div> <div style="width: 1%;"></div> </div> <div>89% 10% .</div>
1	M	552	<div> <div style="width: 87%;"></div> <div style="width: 11%;"></div> <div style="width: 2%;"></div> </div> <div>87% 11% .</div>
2	S	335	<div> <div style="width: 70%;"></div> <div style="width: 9%;"></div> <div style="width: 20%;"></div> </div> <div>70% 9% . 20%</div>
2	T	335	<div> <div style="width: 68%;"></div> <div style="width: 10%;"></div> <div style="width: 22%;"></div> </div> <div>68% 10% . 20%</div>

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 13255 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 [NiFe]-hydrogenase 2 large subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	551	Total	C	N	O	S	0	1	0
			4286	2723	741	802	20			
1	M	551	Total	C	N	O	S	0	1	0
			4287	2724	740	803	20			

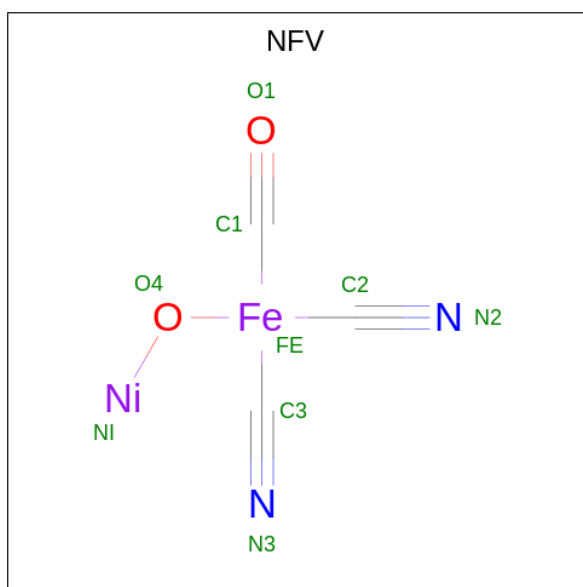
- Molecule 2 is a protein called [NiFe]-hydrogenase 2 small subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	S	268	Total	C	N	O	S	0	1	0
			2046	1297	364	372	13			
2	T	268	Total	C	N	O	S	0	1	0
			2046	1297	364	372	13			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

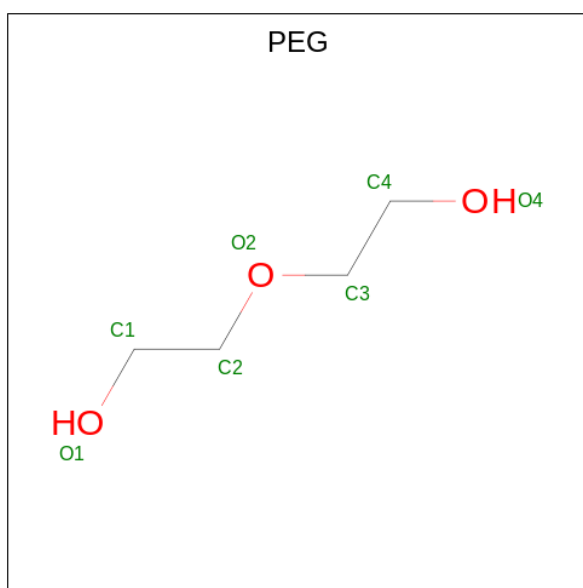
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	1	Total	Mg	0	0
			1	1		
3	M	1	Total	Mg	0	0
			1	1		

- Molecule 4 is NI-FE OXIDIZED ACTIVE CENTER (three-letter code: NFV) (formula: C₃FeN₂NiO₂).



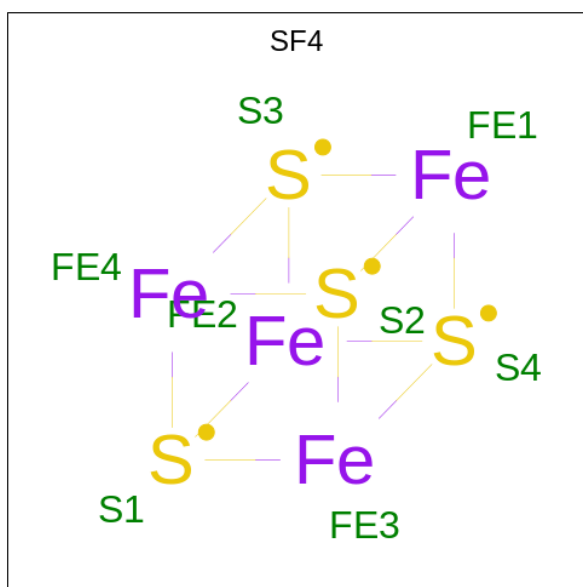
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	L	1	Total	C	Fe	N	Ni	O	0	0
			9	3	1	2	1	2		
4	M	1	Total	C	Fe	N	Ni	O	0	0
			9	3	1	2	1	2		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



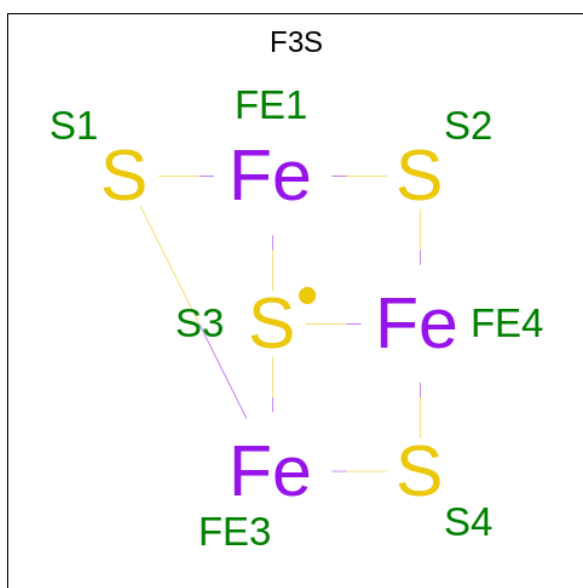
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	L	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe_4S_4).



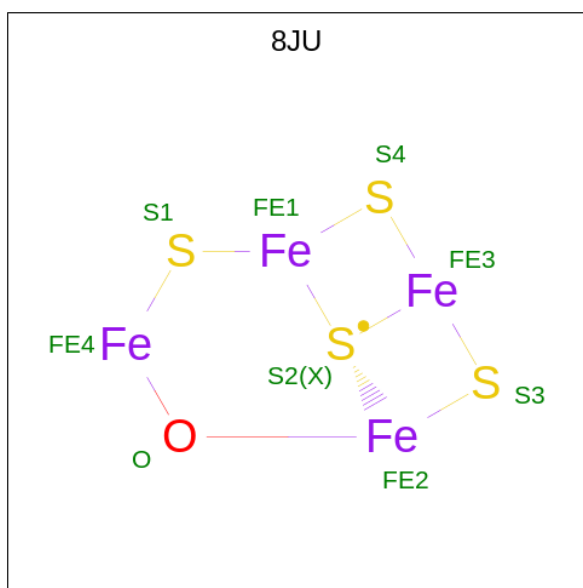
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	S	1	Total	Fe	S	0	0
			8	4	4		
6	T	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 7 is FE3-S4 CLUSTER (three-letter code: F3S) (formula: Fe_3S_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	S	1	Total	Fe	S	0	0
			7	3	4		
7	T	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 8 is FE4-S4-O CLUSTER (three-letter code: 8JU) (formula: Fe_4OS_4).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	S	1	Total	Fe	O	S	0	0
			9	4	1	4		
8	T	1	Total	Fe	O	S	0	0
			9	4	1	4		

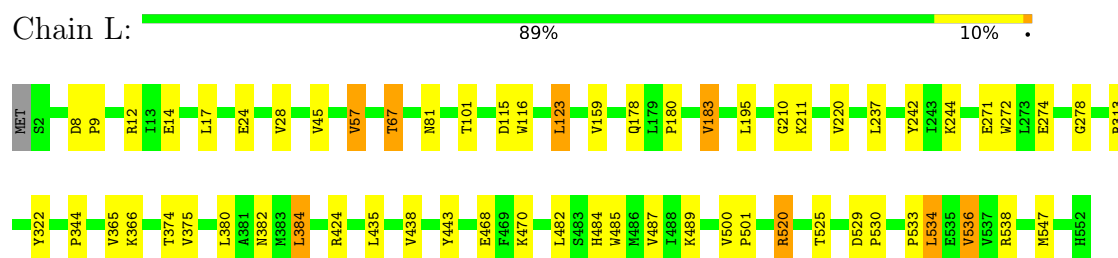
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	L	185	Total	O	0	0
			185	185		
9	S	79	Total	O	0	0
			79	79		
9	M	169	Total	O	0	0
			169	169		
9	T	82	Total	O	0	0
			82	82		

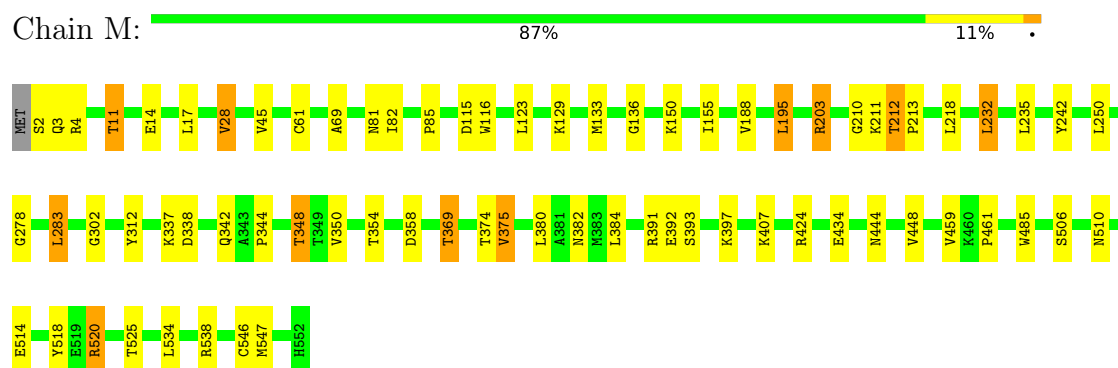
3 Residue-property plots

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.

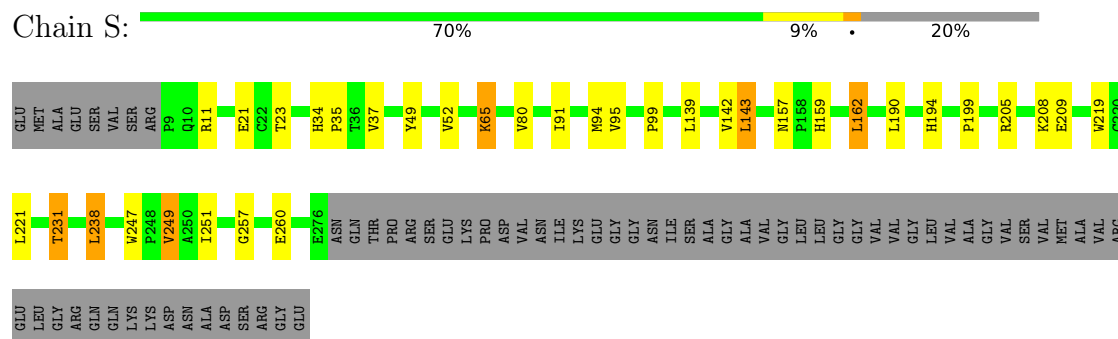
- Molecule 1: [NiFe]-hydrogenase 2 large subunit



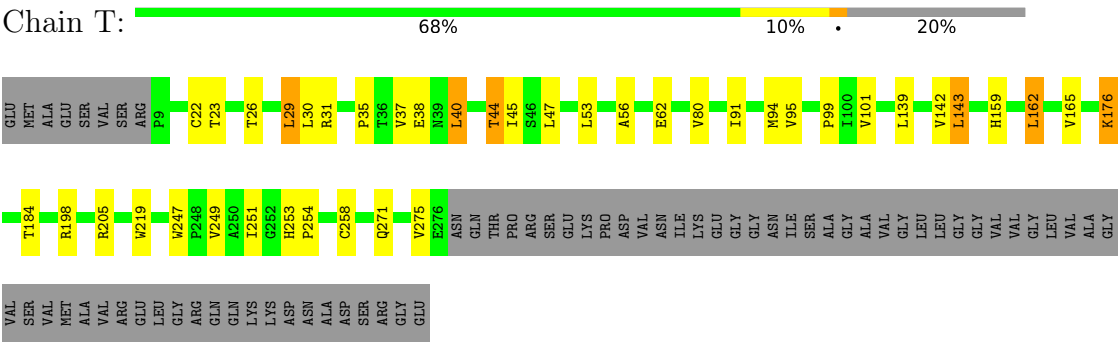
- Molecule 1: [NiFe]-hydrogenase 2 large subunit



- Molecule 2: [NiFe]-hydrogenase 2 small subunit



- Molecule 2: [NiFe]-hydrogenase 2 small subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.97Å 121.83Å 99.17Å 90.00° 102.93° 90.00°	Depositor
Resolution (Å)	37.86 – 2.05 37.86 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.2 (37.86-2.05) 99.3 (37.86-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.37 (at 2.05Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.166 , 0.217 0.166 , 0.217	Depositor DCC
R_{free} test set	5431 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.699	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13255	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% 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: NFV, SF4, CSO, F3S, PEG, 8JU, MG

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	L	0.44	0/4384	0.59	0/5973
1	M	0.44	0/4385	0.62	3/5974 (0.1%)
2	S	0.42	0/2105	0.59	0/2866
2	T	0.42	0/2105	0.60	0/2866
All	All	0.43	0/12979	0.60	3/17679 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	M	232	LEU	CA-CB-CG	8.28	134.34	115.30
1	M	203	ARG	NE-CZ-NH2	-6.06	117.27	120.30
1	M	203	ARG	NE-CZ-NH1	5.73	123.17	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	L	4286	0	4221	31	0
1	M	4287	0	4221	43	0
2	S	2046	0	1969	23	0
2	T	2046	0	1969	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	L	1	0	0	0	0
3	M	1	0	0	0	0
4	L	9	0	0	0	0
4	M	9	0	0	0	0
5	L	7	0	10	3	0
6	S	8	0	0	0	0
6	T	8	0	0	0	0
7	S	7	0	0	0	0
7	T	7	0	0	0	0
8	S	9	0	0	0	0
8	T	9	0	0	0	0
9	L	185	0	0	2	0
9	M	169	0	0	4	0
9	S	79	0	0	1	0
9	T	82	0	0	2	0
All	All	13255	0	12390	113	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 4.

All (113) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S:231:THR:HG21	2:S:257:GLY:HA2	1.58	0.86
2:T:91:ILE:HA	2:T:94:MET:HE3	1.60	0.82
2:S:91:ILE:HA	2:S:94:MET:HE3	1.61	0.81
1:M:338:ASP:OD2	1:M:348:THR:HG21	1.83	0.79
1:M:14:GLU:OE1	1:M:546:CSO:OD	2.05	0.71
2:T:40:LEU:HA	2:T:44:THR:HG23	1.78	0.65
1:M:115:ASP:HB3	1:M:538:ARG:HG3	1.77	0.65
2:T:142:VAL:HG12	2:T:143:LEU:HD13	1.78	0.64
2:S:35:PRO:HD2	2:S:162:LEU:HG	1.80	0.63
1:M:212:THR:HG21	9:M:846:HOH:O	2.00	0.62
1:L:435:LEU:HA	1:L:438:VAL:HG13	1.83	0.60
1:M:283:LEU:HB2	1:M:375:VAL:HG21	1.84	0.60
1:L:210:GLY:CA	2:S:251:ILE:HG12	2.32	0.59
1:L:180:PRO:HD2	1:L:183:VAL:HG13	1.83	0.59
1:M:337:LYS:HB3	1:M:350:VAL:HG22	1.83	0.59
2:S:221:LEU:HD12	2:S:249:VAL:HB	1.84	0.58
1:M:212:THR:HG22	1:M:213:PRO:HA	1.86	0.58
1:L:28:VAL:HG13	1:L:525:THR:HB	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:80:VAL:HB	2:T:139:LEU:HD22	1.87	0.57
2:T:35:PRO:HD2	2:T:162:LEU:HG	1.87	0.57
1:L:278:GLY:O	1:L:520:ARG:HD2	2.05	0.56
1:M:210:GLY:CA	2:T:251:ILE:HG12	2.35	0.56
2:S:231:THR:HB	2:S:260:GLU:HB2	1.87	0.55
1:L:313:ARG:HD3	1:L:322:TYR:CE2	2.42	0.54
1:M:17:LEU:HB2	1:M:547:MET:HG3	1.90	0.54
1:L:57:VAL:HG13	1:L:484:HIS:HE1	1.73	0.53
1:M:369:THR:HG23	9:M:837:HOH:O	2.09	0.53
1:L:468:GLU:OE2	1:L:489:LYS:HE3	2.08	0.53
2:S:142:VAL:HG12	2:S:143:LEU:HD13	1.90	0.52
1:M:11:THR:HG21	2:T:56:ALA:O	2.10	0.52
1:L:380:LEU:HB2	1:L:424:ARG:HA	1.92	0.52
1:M:391:ARG:NH1	1:M:434:GLU:OE1	2.43	0.51
1:M:116:TRP:CZ2	1:M:538:ARG:HG2	2.46	0.51
1:L:533:PRO:HB2	1:L:536:VAL:HG13	1.94	0.50
2:T:26:THR:O	2:T:29:LEU:HB2	2.11	0.50
2:T:91:ILE:HA	2:T:94:MET:CE	2.37	0.50
1:M:203:ARG:NH2	9:M:802:HOH:O	2.34	0.50
1:M:302:GLY:HA3	1:M:312:TYR:HE1	1.77	0.50
2:T:91:ILE:HG22	2:T:94:MET:HE3	1.93	0.49
1:M:2:S:OG	1:M:3:GLN:N	2.46	0.49
1:L:178:GLN:HE22	5:L:603:PEG:H12	1.78	0.48
2:S:80:VAL:HB	2:S:139:LEU:HD22	1.95	0.48
1:L:272:TRP:CD1	5:L:603:PEG:H11	2.49	0.48
1:M:195:LEU:HD23	2:T:38:GLU:HB3	1.96	0.47
1:M:380:LEU:HB2	1:M:424:ARG:HA	1.96	0.47
2:T:94:MET:HE1	2:T:99:PRO:HG3	1.94	0.47
1:L:534:LEU:HD22	1:L:538:ARG:HD3	1.96	0.47
1:M:506:S:HB3	1:M:518:TYR:CE2	2.49	0.47
2:T:30:LEU:CD2	2:T:37:VAL:HG11	2.45	0.47
2:S:205:ARG:HB3	2:S:219:TRP:CD2	2.50	0.47
1:M:82:ILE:HD11	1:M:461:PRO:HB3	1.96	0.46
2:S:91:ILE:HG22	2:S:94:MET:HE3	1.96	0.46
2:S:159:HIS:HA	2:S:162:LEU:HB2	1.97	0.46
1:L:365:VAL:HG12	1:L:366:LYS:O	2.16	0.46
1:M:444:ASN:O	1:M:448:VAL:HG13	2.16	0.46
1:L:8:ASP:HA	1:L:9:PRO:HA	1.81	0.46
1:M:129:LYS:O	1:M:133:MET:HG3	2.16	0.46
2:S:49:TYR:CD2	2:S:65:LYS:HB2	2.52	0.45
1:M:510:ASN:HB2	9:M:895:HOH:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:14:GLU:OE2	2:T:23:THR:OG1	2.30	0.45
1:L:210:GLY:HA3	2:S:251:ILE:HG12	1.97	0.45
1:L:271:GLU:HB2	5:L:603:PEG:H22	1.99	0.45
2:S:247:TRP:CZ2	2:S:249:VAL:HG22	2.52	0.45
1:M:302:GLY:HA3	1:M:312:TYR:CE1	2.52	0.44
2:T:101:VAL:HG23	9:T:642:HOH:O	2.17	0.44
1:M:210:GLY:HA3	2:T:251:ILE:HG12	1.99	0.44
1:L:116:TRP:CZ2	1:L:538:ARG:HG2	2.52	0.44
2:T:40:LEU:HD22	2:T:45:ILE:HG12	1.98	0.44
2:S:199:PRO:HG3	2:T:198:ARG:HG2	2.00	0.44
1:M:4:ARG:NH2	1:M:358:ASP:OD2	2.50	0.44
2:T:253:HIS:CG	2:T:254:PRO:HD2	2.53	0.44
1:M:155:ILE:HD11	1:M:188:VAL:HG11	2.00	0.44
1:L:115:ASP:HB3	1:L:538:ARG:HG3	2.00	0.44
1:M:342:GLN:HE22	1:M:348:THR:HB	1.83	0.44
1:L:244:LYS:HB2	1:L:443:TYR:CZ	2.53	0.43
2:S:247:TRP:CZ3	2:S:249:VAL:HG13	2.53	0.43
1:M:129:LYS:HE3	1:M:129:LYS:HB2	1.85	0.43
1:L:500:VAL:CG1	1:L:501:PRO:HD2	2.49	0.43
2:T:251:ILE:HD12	2:T:251:ILE:C	2.40	0.42
1:M:391:ARG:NH1	1:M:393:SER:HB3	2.35	0.42
1:M:344:PRO:HB2	1:M:485:TRP:CD2	2.54	0.42
2:T:205:ARG:HB3	2:T:219:TRP:CD2	2.55	0.42
1:L:529:ASP:HA	1:L:530:PRO:HD2	1.91	0.42
1:L:17:LEU:HB2	1:L:547:MET:HG3	2.00	0.42
2:S:91:ILE:HA	2:S:94:MET:CE	2.40	0.42
2:T:247:TRP:CH2	2:T:249:VAL:HB	2.55	0.42
2:S:11:ARG:NH2	9:S:602:HOH:O	2.43	0.42
2:S:157:ASN:ND2	2:S:238:LEU:HD13	2.35	0.42
1:M:397:LYS:HA	1:M:397:LYS:HD3	1.81	0.42
1:L:67:THR:HG23	1:L:482:LEU:HD22	2.00	0.42
1:M:344:PRO:HB2	1:M:485:TRP:CG	2.55	0.41
9:L:741:HOH:O	1:M:150:LYS:HE2	2.21	0.41
1:L:12:ARG:HA	2:S:52:VAL:O	2.20	0.41
1:L:14:GLU:OE2	2:S:23:THR:OG1	2.35	0.41
1:L:344:PRO:HB2	1:L:485:TRP:CG	2.55	0.41
2:T:159:HIS:HA	2:T:162:LEU:HB2	2.02	0.41
1:M:28:VAL:CG1	1:M:525:THR:HB	2.50	0.41
1:L:123:LEU:HD21	1:L:159:VAL:HG21	2.02	0.41
1:L:538:ARG:NH2	9:L:716:HOH:O	2.54	0.41
2:T:184:THR:HA	9:T:669:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:136:GLY:O	1:M:407:LYS:NZ	2.54	0.41
1:L:384:LEU:HD12	1:L:384:LEU:HA	1.97	0.41
1:M:278:GLY:O	1:M:520:ARG:HD2	2.20	0.41
2:S:94:MET:HE1	2:S:99:PRO:HG3	2.02	0.40
1:M:61:CYS:HA	2:T:22:CYS:HB2	2.03	0.40
1:M:212:THR:HG22	2:T:31:ARG:HH11	1.86	0.40
1:M:510:ASN:HD21	1:M:514[B]:GLU:CG	2.33	0.40
2:T:271:GLN:NE2	2:T:275:VAL:O	2.44	0.40
1:L:470:LYS:HG2	1:L:487:VAL:HG22	2.04	0.40
2:S:34:HIS:HA	2:S:35:PRO:HA	1.92	0.40
1:M:69:ALA:HB1	1:M:218:LEU:HD12	2.03	0.40
2:T:176:LYS:HB2	2:T:176:LYS:HE2	1.37	0.40
1:M:85:PRO:HG3	1:M:459:VAL:HG23	2.02	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	L	549/552 (100%)	527 (96%)	21 (4%)	1 (0%)	44	38
1	M	549/552 (100%)	527 (96%)	21 (4%)	1 (0%)	44	38
2	S	267/335 (80%)	260 (97%)	7 (3%)	0	100	100
2	T	267/335 (80%)	258 (97%)	9 (3%)	0	100	100
All	All	1632/1774 (92%)	1572 (96%)	58 (4%)	2 (0%)	48	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	211	LYS
1	M	211	LYS

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	L	460/460 (100%)	440 (96%)	20 (4%)	25	19
1	M	460/460 (100%)	438 (95%)	22 (5%)	21	15
2	S	211/261 (81%)	197 (93%)	14 (7%)	14	7
2	T	211/261 (81%)	198 (94%)	13 (6%)	15	8
All	All	1342/1442 (93%)	1273 (95%)	69 (5%)	20	13

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	24	GLU
1	L	45	VAL
1	L	57	VAL
1	L	67	THR
1	L	81	ASN
1	L	101	THR
1	L	123	LEU
1	L	183	VAL
1	L	195	LEU
1	L	220	VAL
1	L	237	LEU
1	L	242	TYR
1	L	274	GLU
1	L	374	THR
1	L	375	VAL
1	L	382	ASN
1	L	384	LEU
1	L	520	ARG
1	L	534	LEU
1	L	536	VAL
2	S	21	GLU
2	S	37	VAL
2	S	65	LYS
2	S	95	VAL

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Mol	Chain	Res	Type
2	S	143	LEU
2	S	162	LEU
2	S	190	LEU
2	S	194	HIS
2	S	208	LYS
2	S	209[A]	GLU
2	S	209[B]	GLU
2	S	231	THR
2	S	238	LEU
2	S	249	VAL
1	M	11	THR
1	M	28	VAL
1	M	45	VAL
1	M	81	ASN
1	M	123	LEU
1	M	195	LEU
1	M	212	THR
1	M	232	LEU
1	M	235	LEU
1	M	242	TYR
1	M	250	LEU
1	M	283	LEU
1	M	348	THR
1	M	354	THR
1	M	369	THR
1	M	374	THR
1	M	375	VAL
1	M	382	ASN
1	M	384	LEU
1	M	392	GLU
1	M	520	ARG
1	M	534	LEU
2	T	29	LEU
2	T	40	LEU
2	T	44	THR
2	T	47	LEU
2	T	53	LEU
2	T	62[A]	GLU
2	T	62[B]	GLU
2	T	95	VAL
2	T	143	LEU
2	T	162	LEU

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Mol	Chain	Res	Type
2	T	165	VAL
2	T	176	LYS
2	T	258	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	CSO	M	546	1,4	3,6,7	0.76	0	0,6,8	-	-
1	CSO	L	546	1,4	3,6,7	0.73	0	0,6,8	-	-

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
1	CSO	M	546	1,4	-	0/1/5/7	-
1	CSO	L	546	1,4	-	0/1/5/7	-

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
1	M	546	CSO	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 2 are monoatomic - leaving 9 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	NFV	L	602	1	3,8,8	2.41	3 (100%)	-		
7	F3S	T	501	2	0,9,9	-	-	-		
4	NFV	M	701	1	3,8,8	3.70	1 (33%)	-		
8	8JU	S	502	2	0,11,11	-	-	-		
8	8JU	T	502	2	0,11,11	-	-	-		
6	SF4	S	500	2	0,12,12	-	-	-		
5	PEG	L	603	-	6,6,6	0.45	0	5,5,5	0.43	0
7	F3S	S	501	2	0,9,9	-	-	-		
6	SF4	T	500	2	0,12,12	-	-	-		

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
7	F3S	T	501	2	-	-	0/3/3/3
8	8JU	S	502	2	-	-	0/2/3/3
8	8JU	T	502	2	-	-	0/2/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	SF4	S	500	2	-	-	0/6/5/5
5	PEG	L	603	-	-	3/4/4/4	-
7	F3S	S	501	2	-	-	0/3/3/3
6	SF4	T	500	2	-	-	0/6/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	M	701	NFV	O1-C1	-6.31	1.06	1.16
4	L	602	NFV	O1-C1	-2.67	1.12	1.16
4	L	602	NFV	C2-N2	-2.34	1.08	1.13
4	L	602	NFV	C3-N3	2.21	1.18	1.13

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	L	603	PEG	O2-C3-C4-O4
5	L	603	PEG	O1-C1-C2-O2
5	L	603	PEG	C1-C2-O2-C3

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	L	603	PEG	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	L	550/552 (99%)	-0.68	0 100 100	4, 11, 26, 43	1 (0%)
1	M	550/552 (99%)	-0.62	0 100 100	5, 13, 25, 47	1 (0%)
2	S	268/335 (80%)	-0.50	0 100 100	6, 14, 33, 51	1 (0%)
2	T	268/335 (80%)	-0.55	0 100 100	6, 14, 28, 50	1 (0%)
All	All	1636/1774 (92%)	-0.61	0 100 100	4, 13, 27, 51	4 (0%)

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CSO	L	546	7/8	0.97	0.08	6,8,10,11	1
1	CSO	M	546	7/8	0.97	0.07	8,10,13,13	1

6.3 Carbohydrates [i](#)

There are no monosaccharides 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(\AA^2)	Q<0.9
5	PEG	L	603	7/7	0.91	0.10	10,20,28,35	0
3	MG	M	700	1/1	0.98	0.02	10,10,10,10	0
6	SF4	S	500	8/8	0.98	0.04	9,13,14,17	0
6	SF4	T	500	8/8	0.98	0.05	10,13,18,18	0
3	MG	L	601	1/1	0.99	0.03	7,7,7,7	0
4	NFV	L	602	9/9	0.99	0.04	4,7,11,12	0
4	NFV	M	701	9/9	0.99	0.04	6,8,11,14	0
7	F3S	S	501	7/7	0.99	0.03	5,6,8,8	0
7	F3S	T	501	7/7	0.99	0.04	7,8,10,12	0
8	8JU	S	502	9/9	0.99	0.04	10,12,19,20	0
8	8JU	T	502	9/9	0.99	0.04	12,14,18,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.