



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:08 PM UTC

PDB ID : 3OYI / pdb_00003oyi
Title : Crystal structure of the PFV S217Q mutant intasome in complex with manganese
Authors : Hare, S.; Cherepanov, P.
Deposited on : 2010-09-23
Resolution : 2.72 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

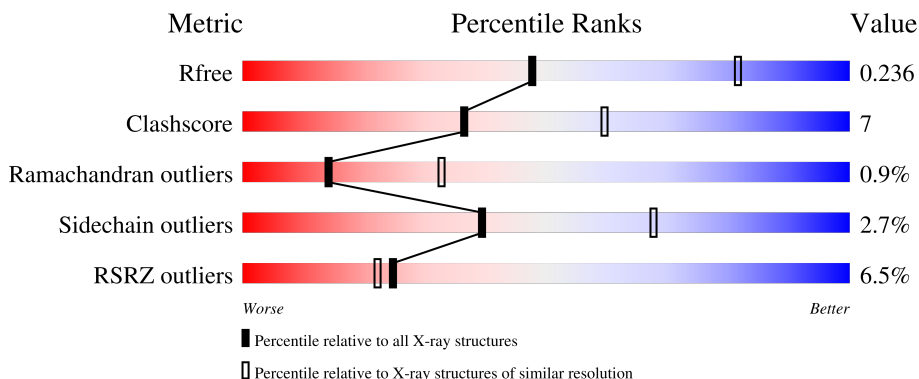
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.72 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	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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	395	 4% 74% 19% • 7%
1	B	395	 5% 39% 6% • 54%
2	C	19	 5% 53% 37% 11%
3	D	17	 6% 53% 35% 12%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 5307 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 PFV integrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	0	0
			2927	1877	514	532	4			
1	B	180	Total	C	N	O	S	0	0	0
			1405	913	228	263	1			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P14350
A	-1	PRO	-	expression tag	UNP P14350
A	0	GLY	-	expression tag	UNP P14350
A	217	GLN	GLY	engineered mutation	UNP P14350
A	218	GLY	SER	variant	UNP P14350
B	-2	GLY	-	expression tag	UNP P14350
B	-1	PRO	-	expression tag	UNP P14350
B	0	GLY	-	expression tag	UNP P14350
B	217	GLN	GLY	engineered mutation	UNP P14350
B	218	GLY	SER	variant	UNP P14350

- Molecule 2 is a DNA chain called DNA (5'-D(*AP*TP*TP*GP*TP*CP*AP*TP*GP*GP*AP*AP*TP*TP*TP*CP*GP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	19	Total	C	N	O	P	0	0	0
			387	187	68	114	18			

- Molecule 3 is a DNA chain called DNA (5'-D(*TP*GP*CP*GP*AP*AP*AP*TP*TP*CP*CP*AP*TP*GP*AP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	17	Total	C	N	O	P	0	0	0
			345	166	65	98	16			

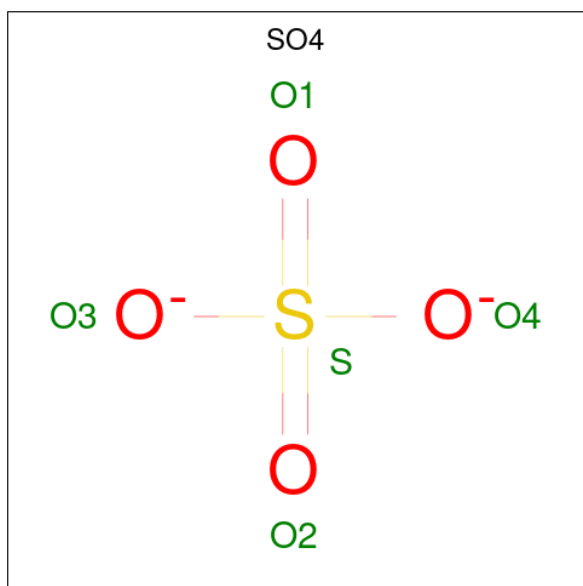
- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

- Molecule 5 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mn	0	0
			1	1		
5	B	1	Total	Mn	0	0
			1	1		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



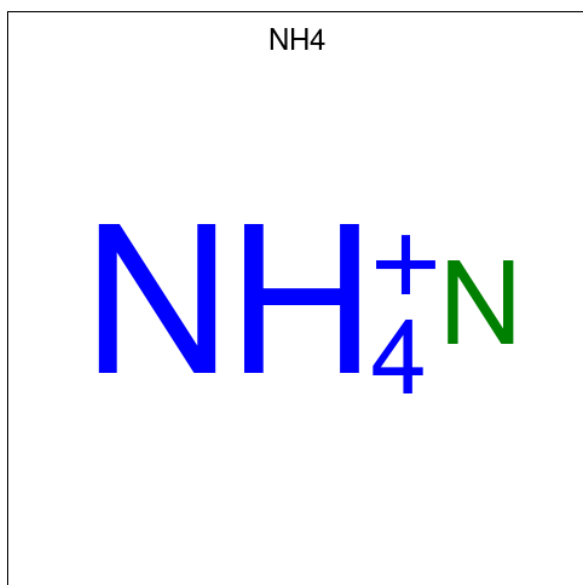
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		

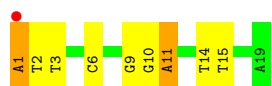
- Molecule 8 is AMMONIUM ION (CCD ID: NH4) (formula: H_4N).



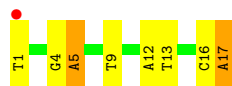
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total N 1 1	0	0

- Molecule 9 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	127	Total O 127 127	0	0
9	B	34	Total O 34 34	0	0
9	C	24	Total O 24 24	0	0
9	D	20	Total O 20 20	0	0



- Molecule 3: DNA (5'-D(*TP*GP*CP*GP*AP*AP*AP*TP*TP*CP*CP*AP*TP*GP*AP*CP*A)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	160.58Å 160.58Å 123.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.95 – 2.72 38.95 – 2.72	Depositor EDS
% Data completeness (in resolution range)	99.1 (38.95-2.72) 99.0 (38.95-2.72)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.73Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.211 , 0.243 0.207 , 0.236	Depositor DCC
R_{free} test set	2188 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	65.2	Xtriage
Anisotropy	0.375	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5307	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, MN, NH4, ZN, GOL

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	0.74	0/3006	0.95	3/4102 (0.1%)
1	B	0.77	0/1442	0.95	0/1971
2	C	0.54	0/433	1.45	4/667 (0.6%)
3	D	0.51	0/387	1.54	8/595 (1.3%)
All	All	0.72	0/5268	1.06	15/7335 (0.2%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1	DT	P-O3'-C3'	7.43	131.35	120.20
3	D	5	DA	C4'-C3'-O3'	-6.68	99.98	110.00
3	D	13	DT	C4'-C3'-O3'	-6.03	100.95	110.00
2	C	14	DT	C4'-C3'-O3'	-5.92	101.11	110.00
2	C	15	DT	C4'-C3'-O3'	-5.89	101.17	110.00
1	A	282	ASP	N-CA-C	5.87	119.26	111.75
3	D	12	DA	C4'-C3'-O3'	-5.86	101.21	110.00
2	C	1	DA	P-O3'-C3'	5.86	128.99	120.20
3	D	9	DT	P-O3'-C3'	5.78	128.87	120.20
2	C	11	DA	C4'-C3'-O3'	-5.65	101.53	110.00
3	D	12	DA	P-O3'-C3'	5.51	128.46	120.20
3	D	17	DA	O4'-C1'-N9	5.38	116.48	108.40
1	A	133	LEU	CA-C-N	-5.30	114.92	120.38
1	A	133	LEU	C-N-CA	-5.30	114.92	120.38
3	D	9	DT	C4'-C3'-O3'	-5.28	102.08	110.00

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	2927	0	2959	48	0
1	B	1405	0	1386	17	0
2	C	387	0	218	7	0
3	D	345	0	192	3	0
4	A	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	5	0	0	0	0
6	B	5	0	0	1	0
7	A	24	0	32	4	0
8	A	1	0	0	0	0
9	A	127	0	0	1	0
9	B	34	0	0	0	0
9	C	24	0	0	0	0
9	D	20	0	0	0	0
All	All	5307	0	4787	71	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 7.

All (71) 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:358:LEU:N	1:A:359:GLY:HA2	1.96	0.81
1:A:358:LEU:H	1:A:359:GLY:HA2	1.54	0.70
1:A:324:GLN:HG2	1:A:340:PRO:HA	1.77	0.67
1:A:142:VAL:HG22	1:A:246:LEU:HD11	1.81	0.62
1:A:276:THR:HG22	1:B:178:ILE:HD12	1.82	0.62
1:A:286:LEU:HB3	1:A:290:GLU:HB2	1.84	0.60
1:A:337:TRP:CG	1:A:370:LYS:HE2	2.38	0.59
1:B:163:THR:HG23	1:B:195:PHE:HB2	1.86	0.58
1:A:287:THR:OG1	1:A:290:GLU:HG3	2.05	0.57
1:A:12:LEU:HD23	1:A:37:VAL:HG21	1.85	0.57
1:A:214:PRO:HA	1:A:217:GLN:NE2	2.20	0.56
1:A:358:LEU:N	1:A:359:GLY:CA	2.68	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:258:SER:O	1:B:261:LEU:O	2.26	0.54
1:B:283:THR:C	1:B:285:ASP:H	2.15	0.54
1:B:172:VAL:O	1:B:175:SER:HB2	2.08	0.54
1:A:118:PRO:HG2	1:A:148:GLY:HA3	1.91	0.53
2:C:1:DA:H2''	2:C:2:DT:H5'	1.91	0.53
1:B:163:THR:HG21	1:B:194:THR:HB	1.90	0.52
1:A:161:PRO:O	1:A:189:ALA:HB2	2.09	0.52
1:A:241:LYS:HG2	7:A:802:GOL:O3	2.09	0.52
1:A:315:TRP:CE2	1:A:371:PRO:HD3	2.44	0.52
1:B:127:ILE:HA	1:B:144:VAL:O	2.08	0.52
1:A:219:LYS:HE3	2:C:6:DC:OP1	2.09	0.52
2:C:1:DA:N3	2:C:2:DT:H71	2.26	0.51
1:A:356:ASP:O	1:A:356:ASP:OD1	2.30	0.50
1:B:264:THR:OG1	1:B:267:GLN:HG3	2.12	0.50
1:A:162:SER:HB2	7:A:801:GOL:H2	1.94	0.49
1:A:76:ILE:HG22	1:A:85:MET:CE	2.43	0.49
1:A:14:GLN:O	1:A:17:GLN:HG2	2.13	0.49
1:A:214:PRO:HA	1:A:217:GLN:HE22	1.77	0.49
2:C:9:DG:H2'	2:C:9:DG:O5'	2.13	0.48
1:B:127:ILE:HG22	1:B:145:VAL:HG22	1.94	0.47
1:A:337:TRP:CD1	1:A:370:LYS:HE2	2.49	0.46
3:D:4:DG:H2''	3:D:5:DA:O5'	2.15	0.46
1:A:91:LYS:HD3	1:A:95:ARG:NH2	2.31	0.46
1:A:325:GLU:HA	1:A:369:LEU:HD23	1.98	0.46
1:A:348:ASN:HB2	1:A:349:PRO:CD	2.47	0.45
1:A:135:PRO:HA	1:A:140:LEU:HD23	1.99	0.45
1:A:156:TYR:OH	1:A:173:LEU:HB2	2.17	0.45
1:A:58:VAL:HG11	1:A:89:VAL:HA	1.98	0.45
1:B:283:THR:C	1:B:285:ASP:N	2.74	0.44
1:A:43:GLU:H	1:A:43:GLU:CD	2.25	0.44
1:B:257:TYR:HA	1:B:264:THR:HG22	1.99	0.44
1:A:127:ILE:HG22	1:A:145:VAL:HG13	2.00	0.44
1:A:221:GLU:O	1:A:224:ASN:HB2	2.18	0.44
1:B:259:PRO:HD2	6:B:394:SO4:O2	2.18	0.44
1:A:250:GLN:NE2	1:B:175:SER:HB3	2.32	0.44
1:A:334:ARG:HD3	9:A:440:HOH:O	2.17	0.44
3:D:16:DC:H2'	3:D:17:DA:H1'	2.01	0.43
1:A:295:LEU:HD21	1:A:299:ARG:NH2	2.33	0.43
1:A:12:LEU:HD23	1:A:37:VAL:CG2	2.49	0.43
1:A:59:LEU:HD12	1:A:59:LEU:HA	1.90	0.42
1:A:133:LEU:O	1:A:134:PRO:C	2.60	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:GLN:HB2	1:A:21:ILE:HD11	2.00	0.42
1:A:54:ARG:O	1:A:58:VAL:HG23	2.19	0.42
1:A:59:LEU:HD13	1:A:92:GLN:OE1	2.19	0.42
1:A:221:GLU:OE1	1:A:221:GLU:HA	2.18	0.42
3:D:16:DC:H5''	3:D:17:DA:OP2	2.20	0.42
1:A:96:CYS:O	1:A:97:GLN:C	2.62	0.41
1:A:115:PRO:HG3	2:C:3:DT:C2	2.55	0.41
1:A:137:GLN:H	7:A:802:GOL:H2	1.85	0.41
1:B:133:LEU:HB3	1:B:134:PRO:HD2	2.02	0.41
1:B:232:THR:HA	1:B:235:LEU:HG	2.03	0.41
1:A:356:ASP:OD1	1:A:358:LEU:HG	2.21	0.41
1:A:304:HIS:CD2	1:A:305:PRO:HD2	2.56	0.41
2:C:10:DG:H2''	2:C:11:DA:C8	2.56	0.41
1:A:183:HIS:CD2	1:A:183:HIS:C	2.98	0.41
7:A:803:GOL:H12	1:B:267:GLN:HB3	2.02	0.40
1:A:170:LEU:O	1:A:174:THR:HG23	2.22	0.40
2:C:10:DG:H2''	2:C:11:DA:H8	1.85	0.40
1:B:244:ASP:O	1:B:247:PRO:HD2	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	366/395 (93%)	347 (95%)	15 (4%)	4 (1%)	11	27
1	B	176/395 (45%)	169 (96%)	6 (3%)	1 (1%)	21	42
All	All	542/790 (69%)	516 (95%)	21 (4%)	5 (1%)	14	33

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	357	HIS
1	A	10	ALA
1	A	69	ARG
1	A	201	GLU
1	B	260	VAL

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	328/354 (93%)	320 (98%)	8 (2%)	43	70
1	B	154/354 (44%)	149 (97%)	5 (3%)	34	62
All	All	482/708 (68%)	469 (97%)	13 (3%)	39	68

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ASP
1	A	36	LYS
1	A	73	LEU
1	A	178	ILE
1	A	286	LEU
1	A	318	VAL
1	A	356	ASP
1	A	364	VAL
1	B	232	THR
1	B	235	LEU
1	B	272	ILE
1	B	282	ASP
1	B	286	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN

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Mol	Chain	Res	Type
1	A	217	GLN
1	A	304	HIS
1	A	374	HIS
1	B	266	HIS

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](#)

Of 10 ligands modelled in this entry, 3 are monoatomic and 1 is modelled with single atom - leaving 6 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	$\# Z > 2$
6	SO4	B	394	-	4,4,4	0.26	0	6,6,6	0.17	0
6	SO4	A	394	-	4,4,4	0.19	0	6,6,6	0.32	0
7	GOL	A	805	-	5,5,5	0.43	0	5,5,5	0.34	0
7	GOL	A	803	-	5,5,5	0.42	0	5,5,5	0.33	0
7	GOL	A	801	-	5,5,5	0.42	0	5,5,5	0.45	0
7	GOL	A	802	-	5,5,5	0.38	0	5,5,5	0.89	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
7	GOL	A	802	-	-	0/4/4/4	-
7	GOL	A	805	-	-	0/4/4/4	-
7	GOL	A	801	-	-	3/4/4/4	-
7	GOL	A	803	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	801	GOL	C1-C2-C3-O3
7	A	801	GOL	O2-C2-C3-O3
7	A	801	GOL	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	394	SO4	1	0
7	A	803	GOL	1	0
7	A	801	GOL	1	0
7	A	802	GOL	2	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	368/395 (93%)	-0.10	15 (4%) 41 38	39, 54, 99, 132	0
1	B	180/395 (45%)	0.54	21 (11%) 9 8	47, 62, 131, 138	0
2	C	19/19 (100%)	-0.30	1 (5%) 32 28	40, 56, 86, 117	0
3	D	17/17 (100%)	-0.61	1 (5%) 28 25	40, 49, 72, 99	0
All	All	584/826 (70%)	0.08	38 (6%) 25 22	39, 57, 117, 138	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	LEU	7.8
1	A	358	LEU	5.3
1	B	298	ILE	4.8
1	B	213	HIS	4.6
1	B	286	LEU	4.6
2	C	1	DA	4.5
1	B	281	GLN	4.5
1	B	288	ARG	4.2
1	B	299	ARG	4.2
1	B	279	ALA	4.2
1	B	284	LEU	3.9
1	B	235	LEU	3.8
1	B	257	TYR	3.4
1	A	10	ALA	3.3
1	A	29	THR	3.2
1	A	357	HIS	3.1
1	B	295	LEU	3.0
1	B	261	LEU	2.9
1	A	15	LEU	2.6
1	B	212	TYR	2.6
1	A	30	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	20	TYR	2.4
1	A	19	HIS	2.4
1	B	296	GLN	2.4
1	B	283	THR	2.3
1	A	40	SER	2.3
1	B	260	VAL	2.3
1	B	119	GLN	2.2
1	B	258	SER	2.2
1	A	359	GLY	2.2
1	A	16	LEU	2.2
1	B	259	PRO	2.1
1	A	14	GLN	2.1
1	B	129	TYR	2.1
3	D	1	DT	2.1
1	B	294	LEU	2.1
1	A	21	ILE	2.1
1	A	24	TYR	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 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
7	GOL	A	801	6/6	0.73	0.30	90,91,92,92	0
6	SO4	B	394	5/5	0.80	0.13	124,124,125,125	0
7	GOL	A	802	6/6	0.81	0.27	88,89,90,90	0
7	GOL	A	805	6/6	0.92	0.17	96,98,99,100	0
6	SO4	A	394	5/5	0.93	0.14	84,86,86,87	0
7	GOL	A	803	6/6	0.94	0.17	70,74,75,76	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	NH4	A	396	1/1	0.96	0.30	34,34,34,34	0
5	MN	A	395	1/1	1.00	0.02	45,45,45,45	0
5	MN	B	393	1/1	1.00	0.01	59,59,59,59	0
4	ZN	A	393	1/1	1.00	0.01	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.