Full wwPDB X-ray Structure Validation Report ⓘ

Dec 6, 2016 – 09:53 AM EST

PDB ID	:	5F0J
Title	:	Structure of retromer VPS26-VPS35 subunits bound to SNX3
Authors	:	Lucas, M.; Gershlick, D.; Vidaurrazaga, A.; Rojas, A.L.; Bonifacino, J.S.; Hierro, A.
Deposited on	:	2015-11-27
Resolution	:	2.70 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 ①) were used in the production of this report:

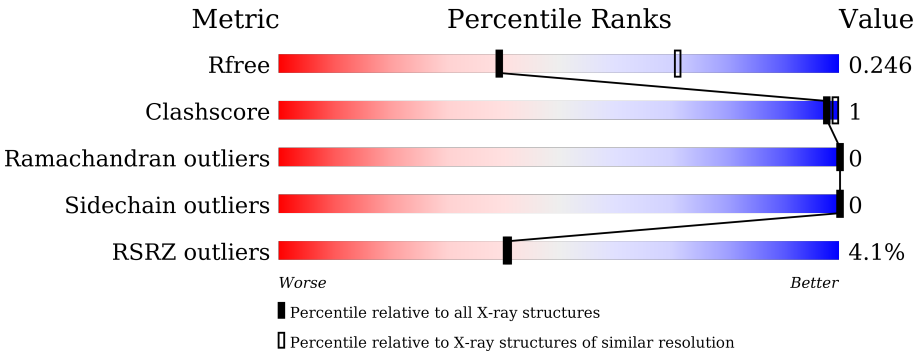
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20028442
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20028442

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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 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 <=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	462	5% 95%
2	B	341	2% 87% 10%
3	C	167	3% 92% 7%

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	C	203	-	-	-	X
5	GOL	B	403	-	-	-	X
5	GOL	B	404	-	-	-	X
6	EDO	A	504	-	-	-	X
6	EDO	A	506	-	-	-	X
6	EDO	B	407	-	-	-	X
6	EDO	B	408	-	-	-	X
6	EDO	B	409	-	-	-	X
6	EDO	B	411	-	-	-	X
6	EDO	C	206	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7732 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 Vacuolar protein sorting-associated protein 35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	458	3709	2359	631	697	22	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	GLY	-	expression tag	UNP Q96QK1
A	10	ALA	-	expression tag	UNP Q96QK1
A	11	MET	-	expression tag	UNP Q96QK1
A	12	GLY	-	expression tag	UNP Q96QK1
A	13	SER	-	expression tag	UNP Q96QK1

- Molecule 2 is a protein called Vacuolar protein sorting-associated protein 26A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
2	B	306	2504	1611	417	466	2	8	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	MSE	-	expression tag	UNP O75436
B	0	GLY	-	expression tag	UNP O75436
B	1	MSE	-	expression tag	UNP O75436
B	327	MSE	-	expression tag	UNP O75436
B	328	GLY	-	expression tag	UNP O75436
B	329	LEU	-	expression tag	UNP O75436
B	330	VAL	-	expression tag	UNP O75436
B	331	PRO	-	expression tag	UNP O75436
B	332	ARG	-	expression tag	UNP O75436
B	333	GLY	-	expression tag	UNP O75436
B	334	SER	-	expression tag	UNP O75436

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Chain	Residue	Modelled	Actual	Comment	Reference
B	335	HIS	-	expression tag	UNP O75436
B	336	HIS	-	expression tag	UNP O75436
B	337	HIS	-	expression tag	UNP O75436
B	338	HIS	-	expression tag	UNP O75436
B	339	HIS	-	expression tag	UNP O75436

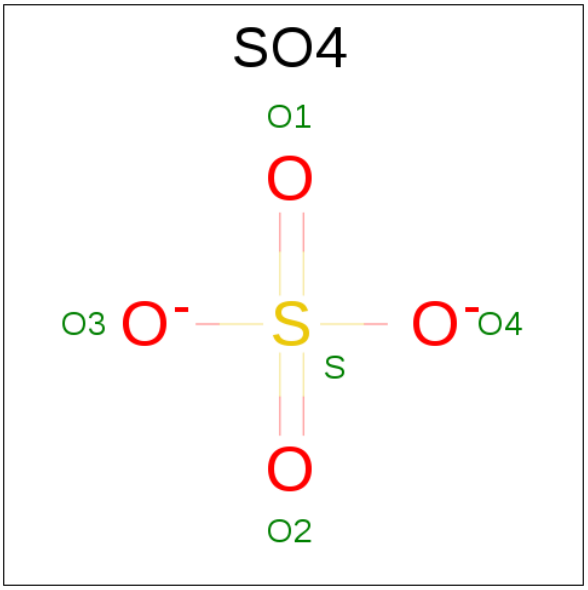
- Molecule 3 is a protein called Sorting nexin-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	155	Total	C	N	O	S	0	0	0
			1268	804	227	235	2			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	GLY	-	expression tag	UNP O60493
C	-3	ALA	-	expression tag	UNP O60493
C	-2	MET	-	expression tag	UNP O60493
C	-1	GLY	-	expression tag	UNP O60493
C	0	SER	-	expression tag	UNP O60493

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

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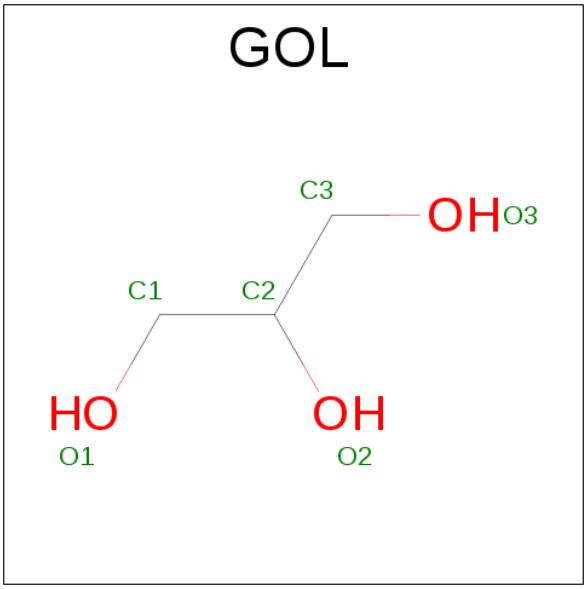
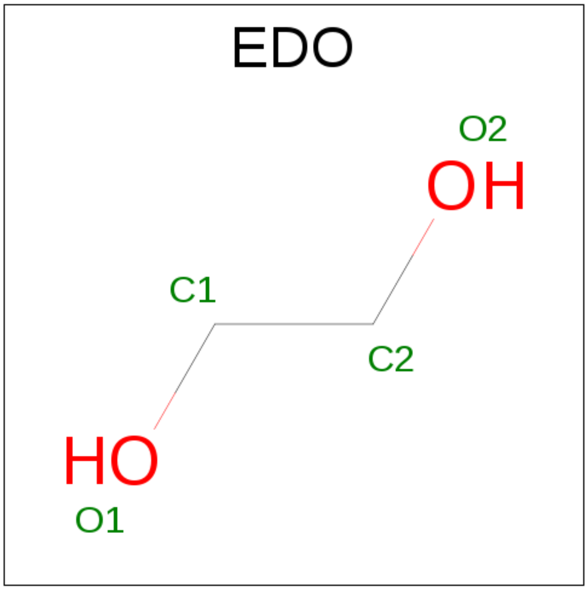
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0

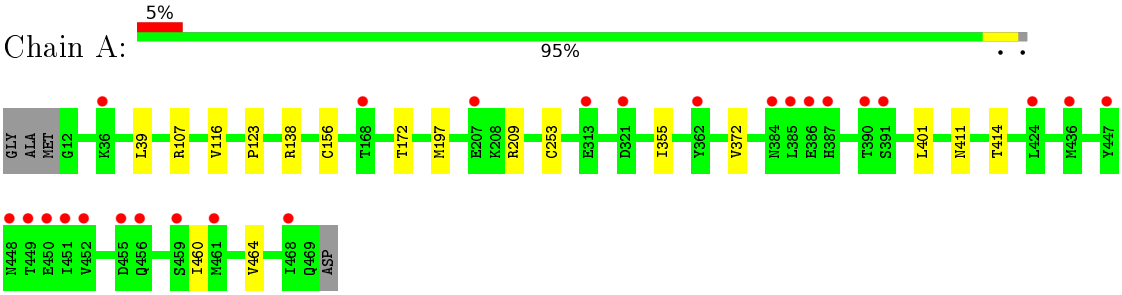
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	29	Total O 29 29	0	0
7	B	47	Total O 47 47	0	0
7	C	19	Total O 19 19	0	0

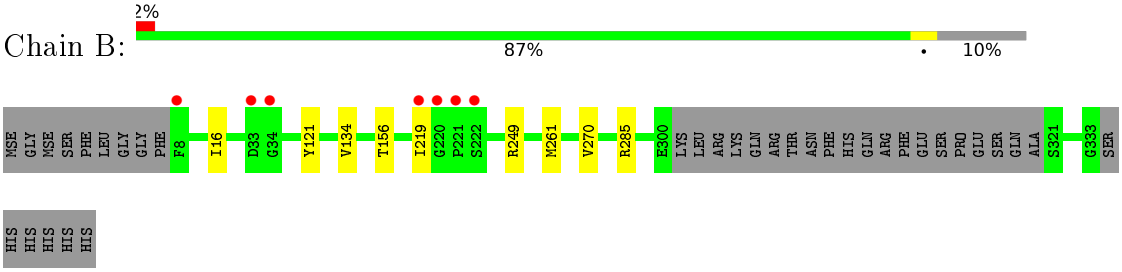
3 Residue-property plots (i)

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.

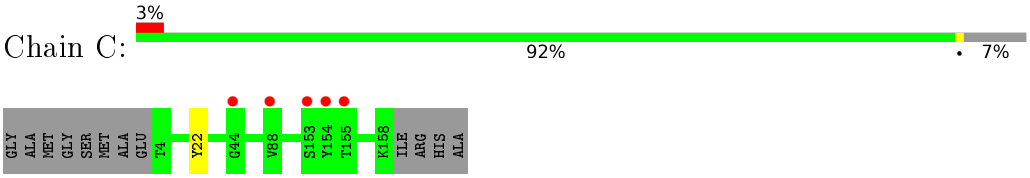
- Molecule 1: Vacuolar protein sorting-associated protein 35



- Molecule 2: Vacuolar protein sorting-associated protein 26A



- Molecule 3: Sorting nexin-3



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	372.82Å 75.58Å 57.53Å 90.00° 97.87° 90.00°	Depositor
Resolution (Å)	56.98 – 2.70 56.99 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.6 (56.98-2.70) 95.7 (56.99-2.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.82 (at 2.69Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.207 , 0.246 0.213 , 0.246	Depositor DCC
R_{free} test set	2099 reflections (5.26%)	DCC
Wilson B-factor (Å^2)	53.3	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 38.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.025 for -h-2*l,-k,l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	7732	wwPDB-VP
Average B, all atoms (Å^2)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.81% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, SO4, EDO

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	0.41	0/3770	0.61	1/5099 (0.0%)
2	B	0.42	0/2546	0.65	1/3417 (0.0%)
3	C	0.43	0/1296	0.65	0/1751
All	All	0.41	0/7612	0.63	2/10267 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	249	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	A	107	ARG	NE-CZ-NH2	5.58	123.09	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	A	3709	0	3762	10	0
2	B	2504	0	2527	7	0
3	C	1268	0	1266	1	0
4	A	10	0	0	0	0
4	B	5	0	0	0	0
4	C	15	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	6	0	8	1	0
5	B	18	0	24	1	0
5	C	6	0	8	0	0
6	A	36	0	54	1	0
6	B	44	0	66	0	0
6	C	16	0	24	0	0
7	A	29	0	0	0	0
7	B	47	0	0	1	0
7	C	19	0	0	0	0
All	All	7732	0	7739	17	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 1.

All (17) 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:355:ILE:HG13	1:A:372:VAL:HG11	1.81	0.62
1:A:460:ILE:O	1:A:464:VAL:HG23	2.14	0.48
1:A:253:CYS:HA	3:C:22:TYR:CG	2.48	0.47
2:B:121:TYR:CD1	2:B:261:MSE:HE1	2.50	0.47
1:A:197:MET:O	1:A:209:ARG:HD2	2.16	0.45
2:B:156:THR:HG22	7:B:523:HOH:O	2.15	0.45
2:B:219:ILE:O	2:B:219:ILE:HD12	2.17	0.44
1:A:355:ILE:HD12	1:A:401:LEU:HG	1.99	0.44
2:B:261:MSE:HE3	2:B:270:VAL:CG2	2.48	0.44
1:A:116:VAL:HG11	1:A:156:CYS:HB3	2.00	0.42
2:B:16:ILE:HD11	2:B:134:VAL:HG23	2.01	0.42
1:A:39:LEU:N	5:A:503:GOL:O2	2.53	0.42
1:A:123:PRO:HB2	1:A:172:THR:HG22	2.01	0.41
2:B:285:ARG:NH1	5:B:403:GOL:O1	2.53	0.41
1:A:411:ASN:O	1:A:414:THR:HG22	2.20	0.41
2:B:261:MSE:CE	2:B:270:VAL:HG21	2.51	0.40
1:A:138:ARG:HA	6:A:506:EDO:H12	2.04	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	456/462 (99%)	448 (98%)	8 (2%)	0	100	100
2	B	302/341 (89%)	295 (98%)	7 (2%)	0	100	100
3	C	153/167 (92%)	149 (97%)	4 (3%)	0	100	100
All	All	911/970 (94%)	892 (98%)	19 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	422/424 (100%)	422 (100%)	0	100	100
2	B	277/298 (93%)	277 (100%)	0	100	100
3	C	141/148 (95%)	141 (100%)	0	100	100
All	All	840/870 (97%)	840 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	141	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

35 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	$\#\|Z\| > 2$
4	SO4	A	501	-	4,4,4	0.39	0	6,6,6	0.09	0
4	SO4	A	502	-	4,4,4	0.37	0	6,6,6	0.15	0
5	GOL	A	503	-	5,5,5	0.30	0	5,5,5	0.16	0
6	EDO	A	504	-	3,3,3	0.59	0	2,2,2	0.31	0
6	EDO	A	505	-	3,3,3	0.46	0	2,2,2	0.20	0
6	EDO	A	506	-	3,3,3	0.34	0	2,2,2	0.49	0
6	EDO	A	507	-	3,3,3	0.50	0	2,2,2	0.09	0
6	EDO	A	508	-	3,3,3	0.37	0	2,2,2	0.63	0
6	EDO	A	509	-	3,3,3	0.48	0	2,2,2	0.26	0
6	EDO	A	510	-	3,3,3	0.51	0	2,2,2	0.26	0
6	EDO	A	511	-	3,3,3	0.41	0	2,2,2	0.54	0
6	EDO	A	512	-	3,3,3	0.54	0	2,2,2	0.12	0
4	SO4	B	401	-	4,4,4	0.33	0	6,6,6	0.29	0
5	GOL	B	402	-	5,5,5	0.17	0	5,5,5	0.31	0
5	GOL	B	403	-	5,5,5	0.31	0	5,5,5	0.22	0
5	GOL	B	404	-	5,5,5	0.46	0	5,5,5	0.48	0
6	EDO	B	405	-	3,3,3	0.48	0	2,2,2	0.33	0
6	EDO	B	406	-	3,3,3	0.40	0	2,2,2	0.52	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
6	EDO	B	407	-	3,3,3	0.45	0	2,2,2	0.38	0
6	EDO	B	408	-	3,3,3	0.47	0	2,2,2	0.24	0
6	EDO	B	409	-	3,3,3	0.50	0	2,2,2	0.32	0
6	EDO	B	410	-	3,3,3	0.43	0	2,2,2	0.22	0
6	EDO	B	411	-	3,3,3	0.50	0	2,2,2	0.31	0
6	EDO	B	412	-	3,3,3	0.47	0	2,2,2	0.30	0
6	EDO	B	413	-	3,3,3	0.46	0	2,2,2	0.38	0
6	EDO	B	414	-	3,3,3	0.62	0	2,2,2	0.31	0
6	EDO	B	415	-	3,3,3	0.51	0	2,2,2	0.23	0
6	EDO	C	201	-	3,3,3	0.56	0	2,2,2	0.09	0
4	SO4	C	202	-	4,4,4	0.26	0	6,6,6	0.16	0
4	SO4	C	203	-	4,4,4	0.26	0	6,6,6	0.35	0
4	SO4	C	204	-	4,4,4	0.37	0	6,6,6	0.21	0
5	GOL	C	205	-	5,5,5	0.22	0	5,5,5	0.25	0
6	EDO	C	206	-	3,3,3	0.53	0	2,2,2	0.08	0
6	EDO	C	207	-	3,3,3	0.44	0	2,2,2	0.43	0
6	EDO	C	208	-	3,3,3	0.57	0	2,2,2	0.07	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	501	-	-	0/0/0/0	0/0/0/0
4	SO4	A	502	-	-	0/0/0/0	0/0/0/0
5	GOL	A	503	-	-	0/4/4/4	0/0/0/0
6	EDO	A	504	-	-	0/1/1/1	0/0/0/0
6	EDO	A	505	-	-	0/1/1/1	0/0/0/0
6	EDO	A	506	-	-	0/1/1/1	0/0/0/0
6	EDO	A	507	-	-	0/1/1/1	0/0/0/0
6	EDO	A	508	-	-	0/1/1/1	0/0/0/0
6	EDO	A	509	-	-	0/1/1/1	0/0/0/0
6	EDO	A	510	-	-	0/1/1/1	0/0/0/0
6	EDO	A	511	-	-	0/1/1/1	0/0/0/0
6	EDO	A	512	-	-	0/1/1/1	0/0/0/0
4	SO4	B	401	-	-	0/0/0/0	0/0/0/0
5	GOL	B	402	-	-	0/4/4/4	0/0/0/0
5	GOL	B	403	-	-	0/4/4/4	0/0/0/0
5	GOL	B	404	-	-	0/4/4/4	0/0/0/0
6	EDO	B	405	-	-	0/1/1/1	0/0/0/0
6	EDO	B	406	-	-	0/1/1/1	0/0/0/0
6	EDO	B	407	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	EDO	B	408	-	-	0/1/1/1	0/0/0/0
6	EDO	B	409	-	-	0/1/1/1	0/0/0/0
6	EDO	B	410	-	-	0/1/1/1	0/0/0/0
6	EDO	B	411	-	-	0/1/1/1	0/0/0/0
6	EDO	B	412	-	-	0/1/1/1	0/0/0/0
6	EDO	B	413	-	-	0/1/1/1	0/0/0/0
6	EDO	B	414	-	-	0/1/1/1	0/0/0/0
6	EDO	B	415	-	-	0/1/1/1	0/0/0/0
6	EDO	C	201	-	-	0/1/1/1	0/0/0/0
4	SO4	C	202	-	-	0/0/0/0	0/0/0/0
4	SO4	C	203	-	-	0/0/0/0	0/0/0/0
4	SO4	C	204	-	-	0/0/0/0	0/0/0/0
5	GOL	C	205	-	-	0/4/4/4	0/0/0/0
6	EDO	C	206	-	-	0/1/1/1	0/0/0/0
6	EDO	C	207	-	-	0/1/1/1	0/0/0/0
6	EDO	C	208	-	-	0/1/1/1	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.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	503	GOL	1	0
6	A	506	EDO	1	0
5	B	403	GOL	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	458/462 (99%)	0.44	25 (5%)	29	27	30, 62, 114, 146	0
2	B	298/341 (87%)	0.21	7 (2%)	64	64	23, 39, 92, 121	0
3	C	155/167 (92%)	0.34	5 (3%)	51	51	28, 58, 105, 116	0
All	All	911/970 (93%)	0.35	37 (4%)	41	41	23, 51, 106, 146	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	451	ILE	6.9
1	A	387	HIS	5.2
2	B	33	ASP	4.9
2	B	221	PRO	4.3
1	A	385	LEU	4.1
2	B	219	ILE	4.0
1	A	386	GLU	3.8
1	A	455	ASP	3.7
1	A	456	GLN	3.7
3	C	88	VAL	3.4
3	C	155	THR	3.4
2	B	34	GLY	3.4
1	A	390	THR	3.4
1	A	362	TYR	3.2
1	A	450	GLU	3.0
1	A	391	SER	3.0
1	A	452	VAL	3.0
2	B	8	PHE	2.8
1	A	384	ASN	2.8
3	C	154	TYR	2.8
3	C	153	SER	2.8
1	A	424	LEU	2.7
1	A	461	MET	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	168	THR	2.7
1	A	448	ASN	2.6
1	A	436	MET	2.6
1	A	207	GLU	2.5
2	B	222	SER	2.5
2	B	220	GLY	2.4
1	A	459	SER	2.3
1	A	468	ILE	2.3
1	A	313	GLU	2.3
1	A	447	TYR	2.3
1	A	36	LYS	2.2
3	C	44	GLY	2.1
1	A	449	THR	2.1
1	A	321	ASP	2.1

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, 95^{th} 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
6	EDO	A	506	4/4	0.94	0.28	4.35	51,52,55,55	0
6	EDO	A	504	4/4	0.87	0.26	4.33	43,44,48,51	0
6	EDO	B	407	4/4	0.90	0.26	3.74	56,56,58,59	0
4	SO4	C	203	5/5	0.82	0.40	3.67	65,67,69,69	5
6	EDO	B	411	4/4	0.81	0.31	3.59	54,54,56,57	0
5	GOL	B	404	6/6	0.78	0.25	3.44	62,66,68,69	0
6	EDO	C	206	4/4	0.85	0.33	3.20	78,79,80,81	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
6	EDO	B	409	4/4	0.91	0.32	2.99	44,46,46,46	0
5	GOL	B	403	6/6	0.82	0.27	2.93	59,65,67,67	0
6	EDO	B	408	4/4	0.84	0.29	2.00	82,83,83,86	0
6	EDO	A	509	4/4	0.81	0.25	1.72	72,73,73,74	0
6	EDO	B	406	4/4	0.93	0.23	1.30	56,60,62,65	0
6	EDO	A	511	4/4	0.95	0.24	1.22	63,64,65,68	0
6	EDO	B	410	4/4	0.94	0.31	0.94	67,67,68,68	0
6	EDO	A	505	4/4	0.85	0.26	0.55	89,93,94,95	0
4	SO4	B	401	5/5	0.96	0.21	0.39	59,59,60,68	0
4	SO4	C	204	5/5	0.95	0.20	0.01	74,75,78,79	0
4	SO4	C	202	5/5	0.99	0.10	-2.00	46,48,49,50	0
6	EDO	B	405	4/4	0.87	0.17	-2.04	47,48,49,51	0
6	EDO	B	415	4/4	0.83	0.26	-	68,70,71,72	0
6	EDO	B	413	4/4	0.91	0.14	-	55,58,60,60	0
6	EDO	A	512	4/4	0.78	0.29	-	65,69,71,72	0
5	GOL	C	205	6/6	0.78	0.27	-	70,75,76,79	0
6	EDO	B	412	4/4	0.86	0.18	-	62,63,64,64	0
6	EDO	A	510	4/4	0.85	0.16	-	70,74,75,77	0
4	SO4	A	502	5/5	0.92	0.14	-	102,102,104,106	0
5	GOL	B	402	6/6	0.91	0.19	-	48,49,51,53	0
6	EDO	C	208	4/4	0.85	0.19	-	51,52,54,55	0
5	GOL	A	503	6/6	0.88	0.16	-	69,74,76,78	0
6	EDO	A	508	4/4	0.90	0.24	-	49,50,50,55	0
6	EDO	A	507	4/4	0.88	0.18	-	70,72,72,72	0
6	EDO	C	201	4/4	0.87	0.26	-	59,64,66,67	0
6	EDO	B	414	4/4	0.83	0.24	-	62,65,65,68	0
6	EDO	C	207	4/4	0.82	0.33	-	84,87,88,91	0
4	SO4	A	501	5/5	0.93	0.16	-	90,91,93,96	0

6.5 Other polymers (i)

There are no such residues in this entry.