



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:38 PM GMT

PDB ID : 1XX1  
Title : Structural basis for ion-coordination and the catalytic mechanism of sphingomyelinases D  
Authors : Murakami, M.T.; Tambourgi, D.V.; Arni, R.K.  
Deposited on : 2004-11-03  
Resolution : 1.75 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

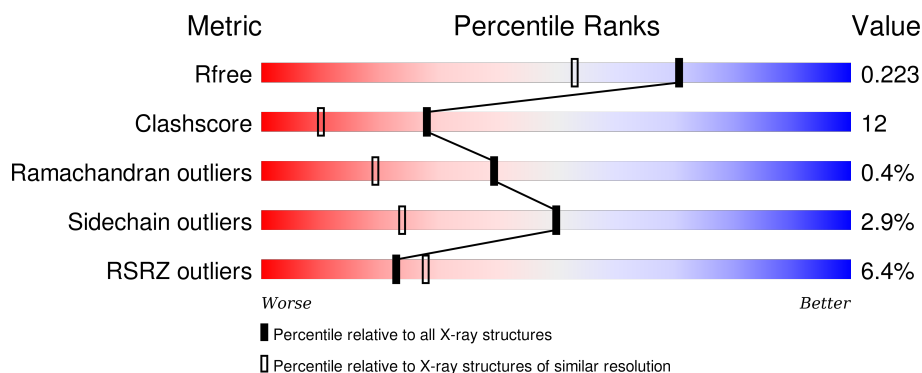
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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	1609 (1.76-1.76)
Clashscore	102246	1730 (1.76-1.76)
Ramachandran outliers	100387	1711 (1.76-1.76)
Sidechain outliers	100360	1711 (1.76-1.76)
RSRZ outliers	91569	1610 (1.76-1.76)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	285	<div> <div>5%</div> <div>82%</div> <div>16%</div> </div>
1	B	285	<div> <div>6%</div> <div>80%</div> <div>19%</div> </div>
1	C	285	<div> <div>8%</div> <div>80%</div> <div>18%</div> </div>
1	D	285	<div> <div>7%</div> <div>80%</div> <div>18%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	4001	-	-	-	X
2	SO4	A	5001	-	-	-	X
2	SO4	B	3002	-	-	X	-
2	SO4	B	4002	-	-	-	X
2	SO4	B	5002	-	-	-	X
2	SO4	C	5003	-	-	-	X
2	SO4	D	5004	-	-	-	X
3	MG	A	9001	-	-	-	X
3	MG	B	9002	-	-	-	X
3	MG	C	9003	-	-	-	X
3	MG	D	9004	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 10258 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 Sphingomyelinase I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	0	0
			2269	1452	382	430	5			
1	B	285	Total	C	N	O	S	0	0	0
			2269	1452	382	430	5			
1	C	285	Total	C	N	O	S	0	0	0
			2269	1452	382	430	5			
1	D	285	Total	C	N	O	S	0	0	0
			2269	1452	382	430	5			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

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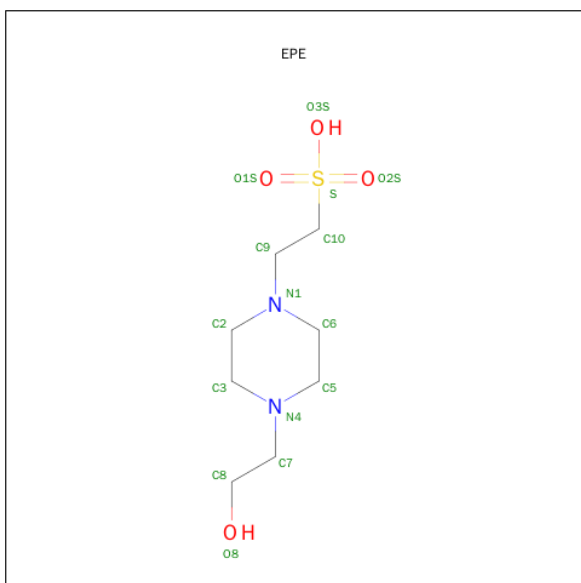
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0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

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C<sub>8</sub>H<sub>18</sub>N<sub>2</sub>O<sub>4</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
4	C	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
4	D	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	273	Total	O	0	0
			273	273		

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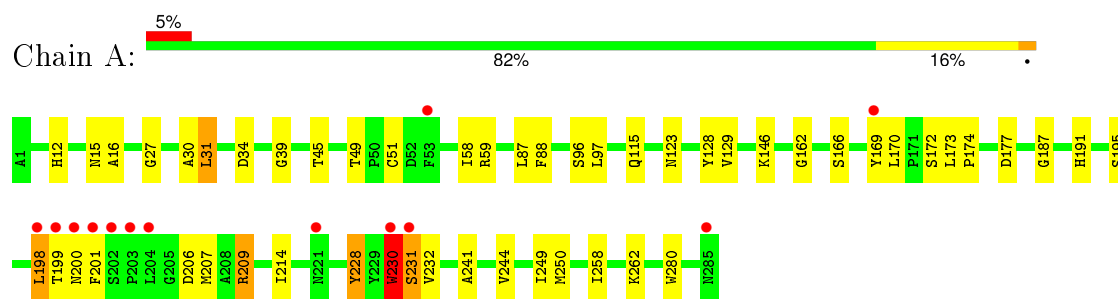
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	283	Total 283	O 283	0	0
5	C	250	Total 250	O 250	0	0
5	D	212	Total 212	O 212	0	0

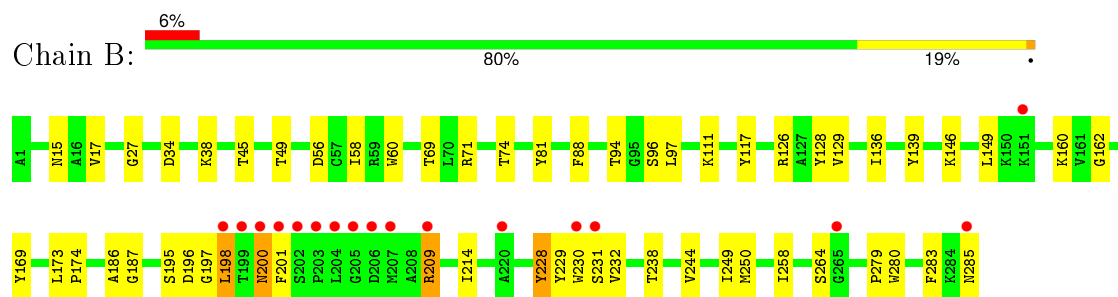
### 3 Residue-property plots

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

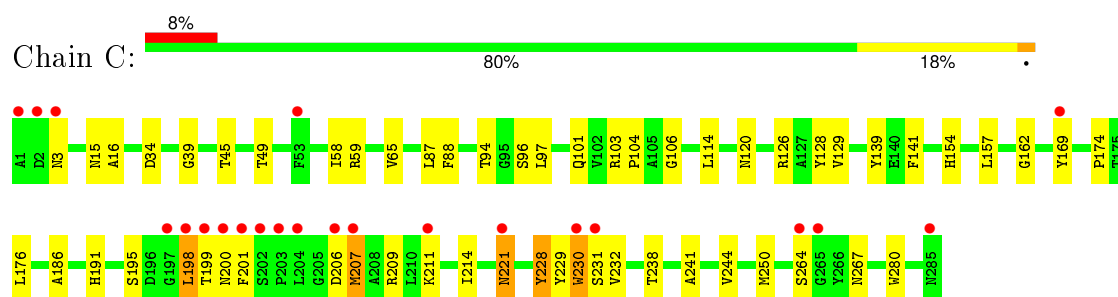
#### • Molecule 1: Spingomyelinase I



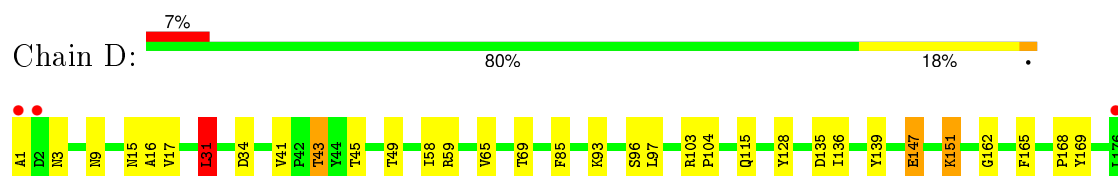
#### • Molecule 1: Spingomyelinase I



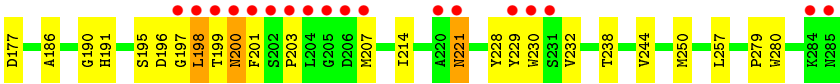
#### • Molecule 1: Spingomyelinase I



#### • Molecule 1: Spingomyelinase I







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	139.82Å 139.82Å 113.46Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 1.75 22.88 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.5 (30.00-1.75) 99.5 (22.88-1.75)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.78 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.186 , 0.225 0.184 , 0.223	Depositor DCC
$R_{free}$ test set	6315 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	20.8	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 30.6	EDS
Estimated twinning fraction	0.082 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 125825 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	10258	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, EPE, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.68	0/2329	0.81	1/3164 (0.0%)
1	B	0.67	0/2329	0.79	0/3164
1	C	0.64	0/2329	0.75	0/3164
1	D	0.59	0/2329	0.74	1/3164 (0.0%)
All	All	0.64	0/9316	0.77	2/12656 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	1
All	All	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	31	LEU	CA-CB-CG	7.85	133.35	115.30
1	A	31	LEU	CA-CB-CG	7.76	133.16	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	230	TRP	Peptide
1	C	230	TRP	Peptide

## 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	2269	0	2190	54	0
1	B	2269	0	2190	63	1
1	C	2269	0	2190	55	1
1	D	2269	0	2190	60	1
2	A	30	0	0	0	0
2	B	30	0	0	4	0
2	C	30	0	0	1	0
2	D	25	0	0	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	15	0	17	0	0
4	C	15	0	17	1	0
4	D	15	0	17	1	0
5	A	273	0	0	10	1
5	B	283	0	0	8	2
5	C	250	0	0	9	1
5	D	212	0	0	5	1
All	All	10258	0	8811	211	4

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 12.

All (211) 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:B:198:LEU:HG	1:B:230:TRP:CZ2	1.63	1.34
1:C:198:LEU:HG	1:C:230:TRP:CZ2	1.75	1.21
1:A:198:LEU:HG	1:A:230:TRP:CZ2	1.74	1.20
1:D:198:LEU:HG	1:D:230:TRP:CZ2	1.80	1.17
1:D:198:LEU:HG	1:D:230:TRP:HZ2	0.96	1.11
1:B:169:TYR:HD2	1:B:173:LEU:HD21	1.17	1.04
1:B:169:TYR:CD2	1:B:173:LEU:HD21	1.93	1.03
1:A:199:THR:HG22	1:B:49:THR:HG21	1.43	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:LEU:CG	1:B:230:TRP:HZ2	1.75	0.98
1:B:198:LEU:HG	1:B:230:TRP:HZ2	0.81	0.97
1:C:198:LEU:HG	1:C:230:TRP:HZ2	1.22	0.97
1:B:111:LYS:HD2	5:B:9249:HOH:O	1.66	0.93
1:D:15:ASN:HD21	1:D:34:ASP:H	1.18	0.90
1:C:15:ASN:HD21	1:C:34:ASP:H	1.19	0.90
1:A:199:THR:HG22	1:B:49:THR:CG2	2.01	0.89
1:A:231:SER:HB2	5:A:9103:HOH:O	1.70	0.88
1:A:15:ASN:HD21	1:A:34:ASP:H	1.17	0.88
1:C:49:THR:OG1	1:D:198:LEU:HD22	1.74	0.87
1:C:3:ASN:HA	5:C:9089:HOH:O	1.74	0.87
1:B:15:ASN:HD21	1:B:34:ASP:H	1.20	0.87
1:A:230:TRP:HB2	1:A:250:MET:HB3	1.56	0.85
1:A:198:LEU:HG	1:A:230:TRP:HZ2	1.42	0.85
2:B:5002:SO4:O1	5:B:9181:HOH:O	1.97	0.82
1:B:56:ASP:HB3	5:B:9188:HOH:O	1.79	0.82
1:A:262:LYS:HE3	5:A:9226:HOH:O	1.79	0.81
1:A:200:ASN:HA	1:A:206:ASP:OD2	1.81	0.81
1:C:198:LEU:HD22	1:D:49:THR:OG1	1.81	0.80
1:B:198:LEU:HD21	2:B:3002:SO4:O2	1.81	0.80
1:B:198:LEU:CD2	2:B:3002:SO4:O2	2.31	0.79
1:A:262:LYS:CE	5:A:9226:HOH:O	2.30	0.79
1:C:230:TRP:HB2	1:C:250:MET:HG2	1.64	0.78
1:B:169:TYR:OH	1:B:198:LEU:HB2	1.84	0.77
1:D:115:GLN:HG3	5:D:9082:HOH:O	1.85	0.76
1:C:97:LEU:HD22	1:C:101:GLN:HB3	1.68	0.75
1:B:126:ARG:HG3	5:B:9263:HOH:O	1.87	0.74
1:D:115:GLN:HG2	5:D:9161:HOH:O	1.86	0.73
1:C:201:PHE:CG	1:D:58:ILE:HD11	2.25	0.70
1:C:230:TRP:CB	1:C:250:MET:HB3	2.21	0.70
1:B:230:TRP:HB2	1:B:250:MET:HB3	1.73	0.69
1:A:49:THR:OG1	1:B:198:LEU:HD22	1.93	0.67
5:C:9140:HOH:O	1:D:97:LEU:CD2	2.43	0.67
1:B:230:TRP:CB	1:B:250:MET:HB3	2.26	0.66
1:C:94:THR:HB	1:C:97:LEU:HD12	1.77	0.66
1:C:199:THR:HG22	1:D:49:THR:CG2	2.25	0.66
1:B:169:TYR:HD2	1:B:173:LEU:CD2	1.99	0.66
1:D:31:LEU:HD13	1:D:85:PHE:CZ	2.30	0.66
1:A:230:TRP:HB2	1:A:250:MET:CB	2.26	0.65
1:A:231:SER:OG	5:A:9104:HOH:O	2.14	0.65
1:C:200:ASN:HA	1:C:206:ASP:OD2	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:ASN:ND2	1:B:34:ASP:H	1.94	0.64
1:D:229:TYR:HD2	1:D:232:VAL:CG2	2.11	0.64
1:A:230:TRP:CB	1:A:250:MET:HB3	2.28	0.64
1:B:174:PRO:HA	5:B:9222:HOH:O	1.97	0.64
1:B:283:PHE:CZ	1:B:285:ASN:HB2	2.33	0.64
1:D:198:LEU:CG	1:D:230:TRP:CZ2	2.70	0.64
1:A:170:LEU:HD22	1:B:60:TRP:CD1	2.32	0.63
1:B:139:TYR:HB2	1:B:186:ALA:HB2	1.80	0.63
1:D:230:TRP:CB	1:D:250:MET:HB3	2.29	0.62
1:A:15:ASN:ND2	1:A:34:ASP:H	1.95	0.62
1:A:166:SER:HA	5:A:9194:HOH:O	1.99	0.62
5:C:9140:HOH:O	1:D:97:LEU:HD21	1.99	0.61
1:B:174:PRO:O	1:B:209:ARG:NH1	2.32	0.61
1:D:15:ASN:ND2	1:D:34:ASP:H	1.97	0.60
1:A:199:THR:CG2	1:B:49:THR:HG21	2.27	0.60
1:C:230:TRP:HB2	1:C:250:MET:CG	2.32	0.59
1:D:41:VAL:O	1:D:43:THR:HG22	2.03	0.59
1:A:58:ILE:HD11	1:B:201:PHE:CG	2.39	0.58
1:A:201:PHE:CG	1:B:58:ILE:HD11	2.38	0.58
1:C:230:TRP:HB2	1:C:250:MET:HB3	1.86	0.58
1:C:15:ASN:ND2	1:C:34:ASP:H	1.98	0.58
1:C:39:GLY:O	1:C:97:LEU:HD21	2.04	0.57
1:A:198:LEU:HD22	1:B:49:THR:OG1	2.03	0.57
1:B:198:LEU:HD23	2:B:3002:SO4:O2	2.04	0.57
1:D:93:LYS:HE3	5:D:9075:HOH:O	2.05	0.57
1:A:49:THR:OG1	1:B:198:LEU:HB3	2.06	0.56
1:D:221:ASN:HD22	1:D:221:ASN:H	1.54	0.56
1:C:195:SER:HB3	1:C:228:TYR:CZ	2.41	0.56
1:C:96:SER:HB3	5:C:9140:HOH:O	2.05	0.56
1:D:229:TYR:HD2	1:D:232:VAL:HG22	1.71	0.55
1:A:232:VAL:HB	1:A:249:ILE:HG23	1.88	0.55
1:C:199:THR:HG22	1:D:49:THR:HG21	1.88	0.55
1:B:231:SER:HB2	5:B:9126:HOH:O	2.05	0.55
1:C:199:THR:HG22	1:D:49:THR:HG23	1.89	0.55
1:A:49:THR:CB	1:B:198:LEU:HD22	2.37	0.55
1:B:88:PHE:CZ	1:B:129:VAL:HG22	2.41	0.55
1:D:203:PRO:O	1:D:207:MET:HG2	2.08	0.54
1:A:169:TYR:OH	1:A:198:LEU:HB3	2.08	0.54
1:D:195:SER:HB3	1:D:228:TYR:CZ	2.42	0.54
1:D:196:ASP:OD1	1:D:197:GLY:N	2.40	0.53
1:C:201:PHE:CB	1:D:58:ILE:HD11	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:135:ASP:OD1	1:D:168:PRO:HD3	2.09	0.53
1:C:65:VAL:HG11	4:C:2003:EPE:H82	1.91	0.52
1:B:195:SER:HB3	1:B:228:TYR:CZ	2.44	0.52
1:B:198:LEU:CG	1:B:230:TRP:CZ2	2.60	0.52
1:C:241:ALA:O	1:C:244:VAL:HG22	2.09	0.52
1:D:17:VAL:HB	1:D:69:THR:HG21	1.91	0.52
1:D:34:ASP:O	1:D:45:THR:HA	2.08	0.52
1:C:162:GLY:HA3	1:C:280:TRP:CH2	2.45	0.51
1:A:262:LYS:HE2	5:A:9226:HOH:O	2.04	0.51
1:C:232:VAL:HG11	1:C:238:THR:HA	1.92	0.51
1:C:34:ASP:O	1:C:45:THR:HA	2.10	0.51
1:B:195:SER:HB3	1:B:228:TYR:CE2	2.45	0.51
1:C:195:SER:HB3	1:C:228:TYR:CE2	2.46	0.50
1:C:88:PHE:CE1	1:C:129:VAL:HG13	2.45	0.50
1:C:221:ASN:HA	5:C:9203:HOH:O	2.11	0.50
1:D:229:TYR:CD2	1:D:232:VAL:HG22	2.46	0.50
1:C:198:LEU:CG	1:C:230:TRP:CZ2	2.69	0.50
1:D:103:ARG:HB3	1:D:104:PRO:HD3	1.94	0.49
1:C:16:ALA:HB2	1:C:59:ARG:HB3	1.94	0.49
1:D:16:ALA:HB2	1:D:59:ARG:HB3	1.93	0.49
1:C:198:LEU:HG	1:C:230:TRP:CH2	2.40	0.49
1:D:230:TRP:HB2	1:D:250:MET:HB3	1.94	0.49
1:D:162:GLY:HA3	1:D:280:TRP:CH2	2.48	0.49
1:B:94:THR:HB	1:B:97:LEU:HD12	1.94	0.49
1:D:195:SER:HB2	1:D:230:TRP:CZ3	2.48	0.48
1:A:174:PRO:HG2	5:A:9194:HOH:O	2.13	0.48
1:C:58:ILE:HD11	1:D:201:PHE:HB2	1.95	0.48
1:C:230:TRP:HB3	1:C:250:MET:HB3	1.94	0.48
1:B:229:TYR:HD2	1:B:232:VAL:CG2	2.26	0.48
1:D:1:ALA:HA	5:D:9180:HOH:O	2.13	0.48
1:D:169:TYR:CE2	1:D:199:THR:HG21	2.48	0.48
1:C:230:TRP:HB2	1:C:250:MET:CB	2.45	0.47
1:A:201:PHE:CD2	1:B:58:ILE:HD11	2.49	0.47
1:D:190:GLY:O	1:D:191:HIS:HB2	2.15	0.47
1:B:34:ASP:O	1:B:45:THR:HA	2.14	0.47
1:B:214:ILE:HD13	1:B:244:VAL:O	2.14	0.47
1:A:230:TRP:HB2	1:A:250:MET:HG2	1.97	0.47
1:D:198:LEU:HD23	2:D:3004:SO4:O1	2.14	0.46
1:A:214:ILE:HD13	1:A:244:VAL:O	2.15	0.46
1:A:241:ALA:O	1:A:244:VAL:HG12	2.15	0.46
1:B:74:THR:O	1:B:126:ARG:HG2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:GLY:O	1:A:97:LEU:HD21	2.16	0.46
1:C:106:GLY:HA3	1:C:141:PHE:O	2.15	0.46
1:A:27:GLY:HA3	1:A:258:ILE:HG13	1.98	0.46
1:C:174:PRO:O	1:C:209:ARG:NH2	2.48	0.46
1:A:88:PHE:CZ	1:A:129:VAL:HG22	2.50	0.46
1:A:195:SER:HB3	1:A:228:TYR:CZ	2.51	0.46
1:B:96:SER:HB3	5:B:9195:HOH:O	2.16	0.46
1:D:214:ILE:HD13	1:D:244:VAL:O	2.16	0.46
1:C:162:GLY:HA2	1:C:191:HIS:HB3	1.99	0.45
1:C:229:TYR:HD2	1:C:232:VAL:CG2	2.29	0.45
1:B:160:LYS:HA	1:B:160:LYS:HD2	1.75	0.45
1:A:199:THR:HG22	1:B:49:THR:HG23	1.89	0.45
1:D:230:TRP:HB3	1:D:250:MET:HB3	1.98	0.45
1:B:17:VAL:HB	1:B:69:THR:HG21	1.98	0.45
1:B:169:TYR:HA	1:B:173:LEU:HD23	1.98	0.45
1:A:199:THR:HG1	1:A:201:PHE:HD1	1.63	0.45
1:B:229:TYR:HD2	1:B:232:VAL:HG23	1.82	0.45
1:D:15:ASN:HD21	1:D:34:ASP:N	1.99	0.44
1:D:195:SER:HB2	1:D:230:TRP:HZ3	1.83	0.44
1:B:15:ASN:HD21	1:B:34:ASP:N	2.01	0.44
1:D:96:SER:HB3	5:D:9167:HOH:O	2.17	0.44
2:C:5003:SO4:O2	5:C:9100:HOH:O	2.21	0.44
1:D:9:ASN:HB2	1:D:257:LEU:HD23	1.99	0.44
1:A:115:GLN:HG3	5:A:9181:HOH:O	2.18	0.44
1:A:34:ASP:O	1:A:45:THR:HA	2.17	0.43
1:D:232:VAL:HG11	1:D:238:THR:HA	2.00	0.43
1:A:195:SER:HA	1:A:228:TYR:O	2.18	0.43
1:B:128:TYR:CG	1:B:279:PRO:HG3	2.53	0.43
1:A:230:TRP:HB2	1:A:250:MET:CG	2.49	0.43
1:A:58:ILE:HD11	1:B:201:PHE:CB	2.49	0.43
1:A:16:ALA:HB2	1:A:59:ARG:HB3	1.99	0.43
1:D:195:SER:HA	1:D:228:TYR:O	2.19	0.43
1:B:232:VAL:HG11	1:B:238:THR:HA	2.01	0.43
1:C:154:HIS:HB3	1:C:157:LEU:HD12	1.99	0.43
1:C:207:MET:HB2	5:C:9242:HOH:O	2.18	0.43
1:A:12:HIS:HB2	1:A:250:MET:HG3	2.01	0.43
1:A:15:ASN:HD21	1:A:34:ASP:N	1.99	0.42
1:A:174:PRO:O	1:A:209:ARG:NH2	2.53	0.42
1:C:201:PHE:HB3	1:D:58:ILE:HD11	2.00	0.42
1:D:65:VAL:HG11	4:D:2004:EPE:H82	2.02	0.42
1:B:195:SER:HB2	1:B:230:TRP:CZ3	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:169:TYR:OH	1:C:198:LEU:HB3	2.19	0.42
1:C:58:ILE:HD11	1:D:201:PHE:CG	2.53	0.42
1:C:103:ARG:HB3	1:C:104:PRO:HD3	2.01	0.42
1:D:147:GLU:O	1:D:151:LYS:HB3	2.20	0.42
1:A:162:GLY:HA3	1:A:280:TRP:CH2	2.55	0.42
1:B:38:LYS:HE3	5:B:9239:HOH:O	2.19	0.42
1:D:162:GLY:HA2	1:D:191:HIS:HB3	2.00	0.42
5:C:9140:HOH:O	1:D:97:LEU:HD23	2.14	0.42
1:C:214:ILE:HD13	1:C:244:VAL:O	2.20	0.42
1:B:196:ASP:O	1:B:229:TYR:HA	2.20	0.42
1:C:126:ARG:HG3	5:C:9094:HOH:O	2.20	0.42
1:D:195:SER:HB3	1:D:228:TYR:CE1	2.55	0.41
1:B:162:GLY:HA3	1:B:280:TRP:CH2	2.55	0.41
1:C:87:LEU:HA	1:C:128:TYR:O	2.20	0.41
1:B:81:TYR:OH	1:B:126:ARG:CZ	2.69	0.41
1:D:169:TYR:CE2	1:D:199:THR:CG2	3.03	0.41
1:C:114:LEU:HG	1:C:120:ASN:ND2	2.35	0.41
1:A:173:LEU:HA	1:A:173:LEU:HD23	1.90	0.41
1:A:51:CYS:HB2	5:A:9102:HOH:O	2.20	0.41
1:D:136:ILE:HG22	1:D:165:PHE:HB3	2.02	0.41
1:D:198:LEU:CB	1:D:230:TRP:CZ2	3.04	0.41
1:C:139:TYR:HB2	1:C:186:ALA:HB2	2.02	0.41
1:B:173:LEU:HA	1:B:173:LEU:HD23	1.60	0.41
1:C:264:SER:HA	1:C:267:ASN:ND2	2.36	0.41
1:A:87:LEU:HA	1:A:128:TYR:O	2.21	0.41
1:D:128:TYR:CG	1:D:279:PRO:HG3	2.56	0.41
1:B:232:VAL:HB	1:B:249:ILE:HG23	2.02	0.41
1:A:30:ALA:HA	1:A:87:LEU:O	2.21	0.41
1:C:176:LEU:HD23	1:C:176:LEU:HA	1.88	0.41
1:C:15:ASN:HD21	1:C:34:ASP:N	2.01	0.40
1:B:71:ARG:HD2	1:B:117:TYR:O	2.20	0.40
1:B:27:GLY:HA3	1:B:258:ILE:HG13	2.03	0.40
1:B:136:ILE:HD13	1:B:174:PRO:HB3	2.02	0.40
1:A:146:LYS:NZ	1:A:187:GLY:O	2.54	0.40
1:A:96:SER:HB3	5:A:9126:HOH:O	2.21	0.40
1:B:146:LYS:NZ	1:B:187:GLY:O	2.54	0.40
1:C:195:SER:HA	1:C:228:TYR:O	2.21	0.40
1:A:162:GLY:HA2	1:A:191:HIS:HB3	2.04	0.40
1:D:139:TYR:HB2	1:D:186:ALA:HB2	2.03	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:200:ASN:O	5:B:9107:HOH:O[4_654]	1.95	0.25
1:B:200:ASN:O	5:C:9128:HOH:O[4_655]	2.00	0.20
1:D:200:ASN:O	5:A:9163:HOH:O[4_654]	2.08	0.12
5:B:9153:HOH:O	5:D:9215:HOH:O[4_655]	2.19	0.01

## 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	283/285 (99%)	278 (98%)	4 (1%)	1 (0%)	39	19
1	B	283/285 (99%)	275 (97%)	6 (2%)	2 (1%)	26	10
1	C	283/285 (99%)	275 (97%)	7 (2%)	1 (0%)	39	19
1	D	283/285 (99%)	275 (97%)	8 (3%)	0	100	100
All	All	1132/1140 (99%)	1103 (97%)	25 (2%)	4 (0%)	39	19

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	231	SER
1	B	198	LEU
1	C	231	SER
1	B	197	GLY

### 5.3.2 Protein sidechains [i](#)

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	240/240 (100%)	231 (96%)	9 (4%)	40	15
1	B	240/240 (100%)	235 (98%)	5 (2%)	61	37
1	C	240/240 (100%)	235 (98%)	5 (2%)	61	37
1	D	240/240 (100%)	231 (96%)	9 (4%)	40	15
All	All	960/960 (100%)	932 (97%)	28 (3%)	50	24

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	LEU
1	A	123	ASN
1	A	172	SER
1	A	177	ASP
1	A	198	LEU
1	A	207	MET
1	A	209	ARG
1	A	228	TYR
1	A	230	TRP
1	B	149	LEU
1	B	200	ASN
1	B	209	ARG
1	B	228	TYR
1	B	264	SER
1	C	198	LEU
1	C	207	MET
1	C	211	LYS
1	C	221	ASN
1	C	228	TYR
1	D	3	ASN
1	D	31	LEU
1	D	43	THR
1	D	147	GLU
1	D	151	LYS
1	D	177	ASP
1	D	198	LEU
1	D	200	ASN
1	D	221	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	ASN
1	A	99	ASN
1	A	122	ASN
1	B	15	ASN
1	B	221	ASN
1	C	15	ASN
1	C	99	ASN
1	C	120	ASN
1	C	123	ASN
1	D	3	ASN
1	D	15	ASN
1	D	115	GLN
1	D	120	ASN
1	D	123	ASN
1	D	200	ASN
1	D	221	ASN
1	D	285	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

Of 30 ligands modelled in this entry, 4 are monoatomic - leaving 26 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	EPE	A	2001	-	14,15,15	0.44	0	18,20,20	1.80	3 (16%)
2	SO4	A	3001	-	4,4,4	0.23	0	6,6,6	0.62	0
2	SO4	A	4001	-	4,4,4	0.42	0	6,6,6	0.28	0
2	SO4	A	5001	-	4,4,4	0.11	0	6,6,6	0.18	0
2	SO4	A	6001	-	4,4,4	0.41	0	6,6,6	0.34	0
2	SO4	A	7001	-	4,4,4	0.29	0	6,6,6	0.35	0
2	SO4	A	8001	-	4,4,4	0.15	0	6,6,6	0.16	0
2	SO4	B	3002	-	4,4,4	0.20	0	6,6,6	0.51	0
2	SO4	B	4002	-	4,4,4	0.30	0	6,6,6	0.45	0
2	SO4	B	5002	-	4,4,4	0.13	0	6,6,6	0.26	0
2	SO4	B	6002	-	4,4,4	0.22	0	6,6,6	0.79	0
2	SO4	B	7002	-	4,4,4	0.31	0	6,6,6	0.24	0
2	SO4	B	8002	-	4,4,4	0.21	0	6,6,6	0.17	0
4	EPE	C	2003	-	14,15,15	0.47	0	18,20,20	1.86	5 (27%)
2	SO4	C	3003	-	4,4,4	0.15	0	6,6,6	0.68	0
2	SO4	C	4003	-	4,4,4	0.31	0	6,6,6	0.25	0
2	SO4	C	5003	-	4,4,4	0.15	0	6,6,6	0.51	0
2	SO4	C	6003	-	4,4,4	0.52	0	6,6,6	0.36	0
2	SO4	C	7003	-	4,4,4	0.22	0	6,6,6	0.13	0
2	SO4	C	8003	-	4,4,4	0.16	0	6,6,6	0.20	0
4	EPE	D	2004	-	14,15,15	0.40	0	18,20,20	2.69	6 (33%)
2	SO4	D	3004	-	4,4,4	0.32	0	6,6,6	0.28	0
2	SO4	D	5004	-	4,4,4	0.25	0	6,6,6	0.20	0
2	SO4	D	6004	-	4,4,4	0.41	0	6,6,6	0.59	0
2	SO4	D	7004	-	4,4,4	0.19	0	6,6,6	0.19	0
2	SO4	D	8004	-	4,4,4	0.07	0	6,6,6	0.15	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	EPE	A	2001	-	-	0/9/19/19	0/1/1/1
2	SO4	A	3001	-	-	0/0/0/0	0/0/0/0
2	SO4	A	4001	-	-	0/0/0/0	0/0/0/0
2	SO4	A	5001	-	-	0/0/0/0	0/0/0/0
2	SO4	A	6001	-	-	0/0/0/0	0/0/0/0
2	SO4	A	7001	-	-	0/0/0/0	0/0/0/0
2	SO4	A	8001	-	-	0/0/0/0	0/0/0/0
2	SO4	B	3002	-	-	0/0/0/0	0/0/0/0
2	SO4	B	4002	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	B	5002	-	-	0/0/0/0	0/0/0/0
2	SO4	B	6002	-	-	0/0/0/0	0/0/0/0
2	SO4	B	7002	-	-	0/0/0/0	0/0/0/0
2	SO4	B	8002	-	-	0/0/0/0	0/0/0/0
4	EPE	C	2003	-	-	0/9/19/19	0/1/1/1
2	SO4	C	3003	-	-	0/0/0/0	0/0/0/0
2	SO4	C	4003	-	-	0/0/0/0	0/0/0/0
2	SO4	C	5003	-	-	0/0/0/0	0/0/0/0
2	SO4	C	6003	-	-	0/0/0/0	0/0/0/0
2	SO4	C	7003	-	-	0/0/0/0	0/0/0/0
2	SO4	C	8003	-	-	0/0/0/0	0/0/0/0
4	EPE	D	2004	-	-	0/9/19/19	0/1/1/1
2	SO4	D	3004	-	-	0/0/0/0	0/0/0/0
2	SO4	D	5004	-	-	0/0/0/0	0/0/0/0
2	SO4	D	6004	-	-	0/0/0/0	0/0/0/0
2	SO4	D	7004	-	-	0/0/0/0	0/0/0/0
2	SO4	D	8004	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	2004	EPE	O2S-S-C10	-5.66	102.07	106.91
4	A	2001	EPE	C9-N1-C6	-3.69	101.80	111.27
4	C	2003	EPE	C3-C2-N1	-3.43	104.49	110.63
4	A	2001	EPE	C5-C6-N1	-3.13	105.03	110.63
4	C	2003	EPE	O1S-S-C10	-3.07	104.29	106.91
4	D	2004	EPE	C5-C6-N1	-2.49	106.18	110.63
4	D	2004	EPE	C3-C2-N1	-2.40	106.34	110.63
4	C	2003	EPE	C5-C6-N1	-2.21	106.67	110.63
4	D	2004	EPE	C7-N4-C5	2.02	116.45	111.27
4	C	2003	EPE	C7-N4-C5	2.32	117.20	111.27
4	A	2001	EPE	C5-N4-C3	4.52	118.69	108.90
4	C	2003	EPE	C5-N4-C3	4.60	118.86	108.90
4	D	2004	EPE	C5-N4-C3	4.88	119.48	108.90
4	D	2004	EPE	O1S-S-C10	7.39	113.21	106.91

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	3002	SO4	3	0
2	B	5002	SO4	1	0
4	C	2003	EPE	1	0
2	C	5003	SO4	1	0
4	D	2004	EPE	1	0
2	D	3004	SO4	1	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 ⓘ

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	285/285 (100%)	0.07	13 (4%)	36 42	11, 18, 35, 59	0
1	B	285/285 (100%)	0.11	17 (5%)	25 30	12, 19, 35, 60	0
1	C	285/285 (100%)	0.24	22 (7%)	16 20	13, 22, 42, 58	0
1	D	285/285 (100%)	0.52	21 (7%)	17 23	13, 25, 49, 72	0
All	All	1140/1140 (100%)	0.23	73 (6%)	23 28	11, 21, 42, 72	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	204	LEU	18.2
1	D	201	PHE	15.8
1	D	198	LEU	12.9
1	B	201	PHE	12.9
1	B	198	LEU	12.4
1	A	198	LEU	12.0
1	D	1	ALA	11.3
1	D	203	PRO	10.8
1	C	201	PHE	10.6
1	C	198	LEU	10.3
1	D	200	ASN	9.6
1	D	199	THR	9.2
1	D	202	SER	8.5
1	C	1	ALA	8.2
1	A	201	PHE	7.8
1	A	199	THR	7.2
1	A	204	LEU	7.2
1	C	2	ASP	7.2
1	B	200	ASN	7.1
1	A	200	ASN	7.1
1	D	205	GLY	6.7

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Mol	Chain	Res	Type	RSRZ
1	A	230	TRP	6.7
1	C	285	ASN	6.6
1	B	199	THR	6.5
1	D	206	ASP	6.4
1	B	230	TRP	6.4
1	B	204	LEU	6.2
1	C	199	THR	6.0
1	D	2	ASP	5.5
1	D	285	ASN	5.2
1	C	230	TRP	4.7
1	C	204	LEU	4.5
1	C	200	ASN	4.1
1	A	285	ASN	4.1
1	C	207	MET	4.1
1	C	203	PRO	4.0
1	C	197	GLY	3.9
1	A	202	SER	3.8
1	D	220	ALA	3.8
1	C	202	SER	3.6
1	A	169	TYR	3.5
1	D	207	MET	3.4
1	D	197	GLY	3.4
1	C	221	ASN	3.4
1	B	202	SER	3.2
1	B	205	GLY	3.1
1	C	231	SER	3.0
1	B	209	ARG	3.0
1	B	206	ASP	2.9
1	B	231	SER	2.9
1	A	53	PHE	2.9
1	B	285	ASN	2.9
1	D	284	LYS	2.7
1	C	3	ASN	2.7
1	D	221	ASN	2.5
1	C	53	PHE	2.5
1	B	203	PRO	2.5
1	A	231	SER	2.4
1	C	265	GLY	2.4
1	D	230	TRP	2.3
1	C	264	SER	2.3
1	D	229	TYR	2.3
1	B	220	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	176	LEU	2.2
1	B	265	GLY	2.2
1	A	221	ASN	2.2
1	A	203	PRO	2.1
1	B	207	MET	2.1
1	C	211	LYS	2.1
1	D	231	SER	2.1
1	B	151	LYS	2.0
1	C	206	ASP	2.0
1	C	169	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 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	MG	A	9001	1/1	1.00	0.16	9.07	7,7,7,7	0
2	SO4	C	5003	5/5	0.92	0.20	7.91	48,48,49,51	0
3	MG	B	9002	1/1	1.00	0.12	7.90	6,6,6,6	0
3	MG	D	9004	1/1	1.00	0.14	7.69	8,8,8,8	0
3	MG	C	9003	1/1	1.00	0.15	6.23	7,7,7,7	0
2	SO4	A	5001	5/5	0.73	0.27	5.86	60,63,63,64	0
2	SO4	B	4002	5/5	0.97	0.14	5.36	29,34,35,36	0
2	SO4	A	4001	5/5	0.95	0.14	5.26	23,30,32,33	0
2	SO4	D	5004	5/5	0.95	0.15	4.91	52,53,54,54	0
2	SO4	B	5002	5/5	0.93	0.15	4.09	45,47,48,48	0
2	SO4	B	6002	5/5	0.97	0.09	0.73	29,31,31,33	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SO4	A	6001	5/5	0.99	0.08	0.71	20,21,23,23	0
4	EPE	A	2001	15/15	0.98	0.08	0.32	17,18,21,23	0
2	SO4	D	6004	5/5	0.99	0.07	-0.09	19,20,20,22	0
2	SO4	B	3002	5/5	0.95	0.10	-0.25	29,34,34,36	0
4	EPE	C	2003	15/15	0.98	0.09	-0.26	17,18,24,28	0
2	SO4	C	3003	5/5	0.94	0.09	-0.29	29,30,31,32	0
2	SO4	A	3001	5/5	0.93	0.09	-0.33	35,35,36,39	0
2	SO4	D	3004	5/5	0.95	0.08	-0.37	28,28,28,30	0
2	SO4	C	4003	5/5	0.96	0.11	-0.88	48,48,49,49	0
2	SO4	C	6003	5/5	0.99	0.06	-1.40	17,18,19,19	0
4	EPE	D	2004	15/15	0.99	0.06	-1.43	18,20,24,27	0
2	SO4	D	8004	5/5	0.92	0.17	-	50,51,51,52	0
2	SO4	A	7001	5/5	0.97	0.09	-	44,45,47,47	0
2	SO4	D	7004	5/5	0.93	0.10	-	63,64,64,64	0
2	SO4	B	7002	5/5	0.96	0.07	-	46,48,49,49	0
2	SO4	C	8003	5/5	0.94	0.16	-	51,51,53,53	0
2	SO4	A	8001	5/5	0.95	0.12	-	46,47,48,49	0
2	SO4	B	8002	5/5	0.94	0.15	-	55,55,56,56	0
2	SO4	C	7003	5/5	0.96	0.12	-	51,52,53,53	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.