



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:19 PM UTC

PDB ID : 6MOE / pdb_00006moe
Title : Monomeric DARPin E2 complex with EpoR
Authors : Jude, K.M.; Mohan, K.; Garcia, K.C.
Deposited on : 2018-10-04
Resolution : 2.09 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

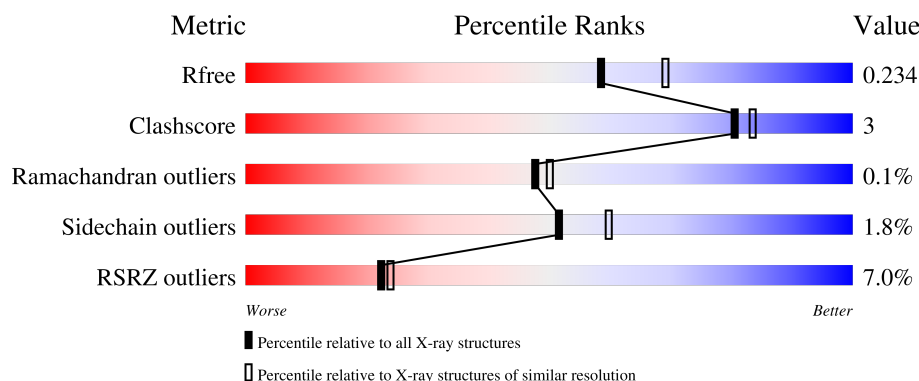
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	168	% 91% 5% .
1	B	168	3% 92% . . .
2	C	229	7% 88% 6% . 5%
2	D	229	13% 85% 8% . 6%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6249 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 E2 DARPin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	161	Total	C	N	O	S	0	0	0
			1194	751	203	237	3			
1	B	161	Total	C	N	O	S	0	0	0
			1194	751	203	237	3			

- Molecule 2 is a protein called Erythropoietin receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	217	Total	C	N	O	S	0	0	0
			1676	1064	293	312	7			
2	D	216	Total	C	N	O	S	0	0	0
			1669	1061	290	311	7			

There are 26 discrepancies between the modelled and reference sequences:

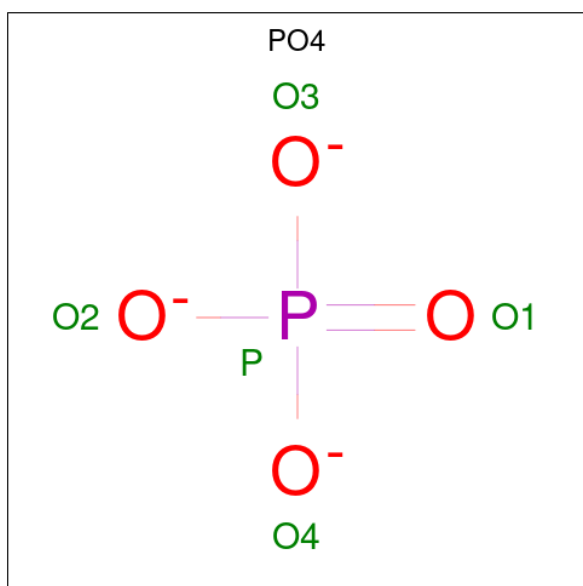
Chain	Residue	Modelled	Actual	Comment	Reference
C	3	PHE	-	expression tag	UNP P19235
C	4	ALA	-	expression tag	UNP P19235
C	5	GLY	-	expression tag	UNP P19235
C	6	SER	-	expression tag	UNP P19235
C	7	ALA	-	expression tag	UNP P19235
C	52	GLN	ASN	conflict	UNP P19235
C	164	GLN	ASN	conflict	UNP P19235
C	226	LYS	-	expression tag	UNP P19235
C	227	GLU	-	expression tag	UNP P19235
C	228	LYS	-	expression tag	UNP P19235
C	229	ALA	-	expression tag	UNP P19235
C	230	ALA	-	expression tag	UNP P19235
C	231	ALA	-	expression tag	UNP P19235
D	3	PHE	-	expression tag	UNP P19235
D	4	ALA	-	expression tag	UNP P19235
D	5	GLY	-	expression tag	UNP P19235

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Chain	Residue	Modelled	Actual	Comment	Reference
D	6	SER	-	expression tag	UNP P19235
D	7	ALA	-	expression tag	UNP P19235
D	52	GLN	ASN	conflict	UNP P19235
D	164	GLN	ASN	conflict	UNP P19235
D	226	LYS	-	expression tag	UNP P19235
D	227	GLU	-	expression tag	UNP P19235
D	228	LYS	-	expression tag	UNP P19235
D	229	ALA	-	expression tag	UNP P19235
D	230	ALA	-	expression tag	UNP P19235
D	231	ALA	-	expression tag	UNP P19235

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

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

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).

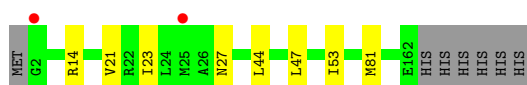


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	104	Total	O	0	0
			104	104		
6	B	101	Total	O	0	0
			101	101		
6	C	116	Total	O	0	0
			116	116		
6	D	127	Total	O	0	0
			127	127		

- Molecule 1: E2 DARPIn



-
- | Category | Number of Genes |
|----------|-----------------|
| MET | 1 |
| G2 | 2 |
| S3 | 2 |
| R14 | 1 |
| L44 | 1 |
| E85 | 1 |
| E88 | 2 |
| Y100 | 1 |
| I119 | 1 |
| V122 | 1 |
| K125 | 1 |
| V130 | 2 |
| L161 | 2 |
| E162 | 1 |
| HIS | 1 |

- | PHE | ALA |
|------|-----|
| A7 | ALA |
| D8 | ALA |
| P9 | ALA |
| K10 | ALA |
| R32 | |
| L33 | |
| E34 | |
| A46 | |
| G47 | |
| V48 | |
| D58 | |
| L59 | |
| E60 | |
| W64 | |
| C63 | |
| T57 | |
| R108 | |
| E157 | |
| V158 | |
| D159 | |
| A162 | |
| G163 | |
| Q164 | |
| G165 | |
| A166 | |
| I167 | |
| S168 | |
| V169 | |
| Q170 | |
| R197 | |
| A198 | |
| R199 | |
| F205 | |
| G206 | |
| G207 | |
| F208 | |
| W212 | |
| D223 | |
| LEU | |
| ASP | |
| LYS | |
| GLU | |
| VAL | |

-
- | Protein | Number of Mutations |
|---------|---------------------|
| PHE | 1 |
| ALA | 1 |
| GLY | 1 |
| SER | 1 |
| ALA | 1 |
| D8 | 1 |
| P9 | 1 |
| A10 | 1 |
| F11 | 1 |
| R21 | 1 |
| R32 | 1 |
| L33 | 1 |
| E34 | 1 |
| D35 | 1 |
| L36 | 1 |
| V37 | 1 |
| A46 | 1 |
| G47 | 1 |
| V48 | 1 |
| G49 | 1 |
| P50 | 1 |
| Q58 | 1 |
| L59 | 1 |
| E60 | 1 |
| W64 | 1 |
| Q71 | 1 |
| C83 | 1 |
| S84 | 1 |
| L85 | 1 |
| S91 | 1 |
| S92 | 1 |
| F93 | 1 |
| A103 | 1 |
| T151 | 1 |
| I154 | 1 |
| A162 | 1 |
| G163 | 1 |
| Q164 | 1 |
| G165 | 1 |
| A166 | 1 |
| G167 | 1 |
| S168 | 1 |
| F169 | 1 |
| I174 | 1 |
| L175 | 1 |
| P176 | 1 |

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.25Å 122.16Å 132.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.77 – 2.09 29.77 – 2.09	Depositor EDS
% Data completeness (in resolution range)	99.5 (29.77-2.09) 84.3 (29.77-2.09)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.30 (at 2.10Å)	Xtriage
Refinement program	PHENIX (1.14_3211: ???)	Depositor
R, R_{free}	0.204 , 0.230 0.206 , 0.234	Depositor DCC
R_{free} test set	2001 reflections (2.98%)	wwPDB-VP
Wilson B-factor (Å ²)	43.0	Xtriage
Anisotropy	0.491	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 53.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6249	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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: PO4, EPE, 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	0.11	0/1209	0.29	0/1644
1	B	0.11	0/1209	0.30	0/1644
2	C	0.14	0/1722	0.38	0/2352
2	D	0.14	0/1715	0.44	2/2342 (0.1%)
All	All	0.13	0/5855	0.37	2/7982 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	202	GLU	CA-C-N	5.32	139.77	127.00
2	D	202	GLU	C-N-CA	5.32	139.77	127.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1194	0	1194	4	0
1	B	1194	0	1194	5	0
2	C	1676	0	1620	10	0
2	D	1669	0	1615	11	0
3	A	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	5	0	0	0	0
3	C	15	0	0	0	0
3	D	15	0	0	0	0
4	A	4	0	6	0	0
4	C	4	0	6	0	0
5	B	15	0	17	1	0
6	A	104	0	0	0	0
6	B	101	0	0	1	0
6	C	116	0	0	1	0
6	D	127	0	0	0	0
All	All	6249	0	5652	30	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 3.

All (30) 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:14:ARG:HB2	1:B:44:LEU:HD13	1.75	0.68
1:A:14:ARG:HB2	1:A:44:LEU:HD13	1.77	0.66
2:C:199:ARG:NH2	2:C:207:GLY:O	2.33	0.62
2:C:223:ASP:OD1	2:C:223:ASP:N	2.21	0.60
2:D:58:GLN:NE2	2:D:60:GLU:O	2.33	0.58
2:C:197:ARG:HG3	2:C:212:TRP:CE3	2.38	0.58
1:A:23:ILE:O	1:A:27:ASN:ND2	2.33	0.57
2:D:197:ARG:HD2	2:D:209:TRP:HB3	1.87	0.56
2:D:162:ALA:HB3	2:D:166:ALA:HB3	1.89	0.54
2:C:32:ARG:HB2	2:C:34:GLU:HG2	1.95	0.48
2:D:58:GLN:HB2	2:D:64:TRP:CD2	2.48	0.48
1:A:47:LEU:HD12	1:A:81:MET:HG3	1.96	0.48
2:D:10:LYS:NZ	2:D:32:ARG:HH22	2.12	0.47
2:D:58:GLN:HB2	2:D:64:TRP:CE2	2.50	0.46
2:C:58:GLN:HB2	2:C:64:TRP:CD2	2.51	0.46
2:D:11:PHE:CE2	2:D:37:VAL:HG11	2.51	0.46
2:C:157:GLU:OE1	2:C:197:ARG:NH2	2.27	0.45
2:C:170:GLN:NE2	6:C:403:HOH:O	2.33	0.45
1:B:100:TYR:CE2	5:B:202:EPE:H21	2.53	0.44
1:B:85:GLU:H	1:B:85:GLU:CD	2.23	0.43
1:A:21:VAL:HG11	1:A:53:ILE:HG23	1.99	0.43
2:C:58:GLN:HB2	2:C:64:TRP:CE2	2.54	0.42
1:B:125:LYS:NZ	6:B:303:HOH:O	2.47	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:50:PRO:HD3	2:D:71:GLN:OE1	2.20	0.42
2:C:159:ASP:HB2	2:C:212:TRP:CH2	2.55	0.42
1:B:119:ILE:HA	1:B:122:VAL:HG12	2.02	0.42
2:D:175:LEU:HD12	2:D:178:ARG:NH2	2.36	0.41
2:D:174:ILE:HD12	2:D:181:CYS:HB3	2.03	0.40
2:C:164:GLN:OE1	2:C:164:GLN:N	2.50	0.40
2:D:151:THR:HA	2:D:154:ILE:HD12	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	A	159/168 (95%)	159 (100%)	0	0	100	100
1	B	159/168 (95%)	159 (100%)	0	0	100	100
2	C	215/229 (94%)	208 (97%)	7 (3%)	0	100	100
2	D	214/229 (93%)	207 (97%)	6 (3%)	1 (0%)	24	22
All	All	747/794 (94%)	733 (98%)	13 (2%)	1 (0%)	48	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	164	GLN

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	122/129 (95%)	122 (100%)	0	100	100
1	B	122/129 (95%)	121 (99%)	1 (1%)	73	81
2	C	176/185 (95%)	170 (97%)	6 (3%)	32	35
2	D	176/185 (95%)	172 (98%)	4 (2%)	44	51
All	All	596/628 (95%)	585 (98%)	11 (2%)	51	60

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	85	GLU
2	C	58	GLN
2	C	83	CYS
2	C	108	ARG
2	C	164	GLN
2	C	208	PHE
2	C	223	ASP
2	D	58	GLN
2	D	83	CYS
2	D	164	GLN
2	D	202	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	B	116	HIS
2	C	70	HIS
2	D	70	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

12 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	$\# Z > 2$
5	EPE	B	202	-	15,15,15	0.83	1 (6%)	19,20,20	1.84	3 (15%)
3	PO4	C	301	-	4,4,4	0.96	0	6,6,6	0.47	0
3	PO4	B	201	-	4,4,4	0.99	0	6,6,6	0.45	0
3	PO4	A	202	-	4,4,4	0.95	0	6,6,6	0.48	0
3	PO4	D	303	-	4,4,4	0.94	0	6,6,6	0.49	0
4	EDO	A	203	-	3,3,3	0.42	0	2,2,2	0.35	0
3	PO4	C	302	-	4,4,4	0.97	0	6,6,6	0.46	0
4	EDO	C	304	-	3,3,3	0.42	0	2,2,2	0.39	0
3	PO4	A	201	-	4,4,4	0.94	0	6,6,6	0.44	0
3	PO4	D	301	-	4,4,4	0.96	0	6,6,6	0.46	0
3	PO4	C	303	-	4,4,4	0.95	0	6,6,6	0.46	0
3	PO4	D	302	-	4,4,4	0.98	0	6,6,6	0.46	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	EDO	A	203	-	-	0/1/1/1	-
5	EPE	B	202	-	-	8/9/19/19	0/1/1/1
4	EDO	C	304	-	-	0/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	202	EPE	C10-S	2.74	1.81	1.77

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	202	EPE	C5-N4-C3	4.52	118.58	108.84
5	B	202	EPE	C7-N4-C5	3.49	120.54	111.24
5	B	202	EPE	C7-N4-C3	3.39	120.27	111.24

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	202	EPE	S-C10-C9-N1
5	B	202	EPE	C9-C10-S-O2S
5	B	202	EPE	C9-C10-S-O3S
5	B	202	EPE	N4-C7-C8-O8
5	B	202	EPE	C9-C10-S-O1S
5	B	202	EPE	C10-C9-N1-C2
5	B	202	EPE	C10-C9-N1-C6
5	B	202	EPE	C8-C7-N4-C5

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	202	EPE	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 [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	161/168 (95%)	0.21	2 (1%) 76 78	41, 58, 84, 129	0
1	B	161/168 (95%)	0.29	5 (3%) 51 54	44, 61, 83, 144	0
2	C	217/229 (94%)	0.64	16 (7%) 20 22	46, 66, 109, 171	0
2	D	216/229 (94%)	0.75	30 (13%) 6 6	44, 67, 119, 158	0
All	All	755/794 (95%)	0.51	53 (7%) 22 24	41, 63, 109, 171	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	166	ALA	5.4
2	C	48	VAL	4.1
2	D	203	PRO	4.0
2	C	46	ALA	4.0
2	D	48	VAL	3.8
2	D	11	PHE	3.8
2	C	7	ALA	3.7
2	C	165	GLY	3.5
2	D	46	ALA	3.4
2	C	162	ALA	3.3
2	C	163	GLY	2.9
1	A	2	GLY	2.9
2	D	205	PHE	2.9
2	D	206	GLY	2.9
2	D	162	ALA	2.8
1	B	2	GLY	2.8
2	D	207	GLY	2.8
2	D	166	ALA	2.8
1	A	25	MET	2.7
2	D	223	ASP	2.7
2	C	205	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
2	D	21	ARG	2.7
2	C	167	GLY	2.7
2	D	50	PRO	2.7
2	C	168	SER	2.6
2	D	9	PRO	2.6
2	D	169	VAL	2.4
2	C	87	THR	2.4
2	D	201	ALA	2.4
2	D	103	ALA	2.4
1	B	161	LEU	2.4
2	D	218	LEU	2.3
1	B	88	GLU	2.3
2	D	10	LYS	2.3
2	D	204	SER	2.3
2	C	8	ASP	2.3
2	D	93	PHE	2.3
2	C	169	VAL	2.3
2	D	91	SER	2.2
1	B	130	VAL	2.2
2	D	202	GLU	2.2
2	D	33	LEU	2.2
2	D	49	GLY	2.1
2	D	85	LEU	2.1
2	C	10	LYS	2.1
2	D	59	LEU	2.1
2	D	175	LEU	2.1
2	D	167	GLY	2.1
2	D	35	ASP	2.1
2	D	178	ARG	2.1
1	B	3	SER	2.1
2	C	9	PRO	2.0
2	C	60	GLU	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
5	EPE	B	202	15/15	0.54	0.22	140,154,186,188	0
4	EDO	C	304	4/4	0.59	0.21	99,99,100,100	0
3	PO4	D	303	5/5	0.67	0.28	165,165,165,166	0
3	PO4	C	302	5/5	0.77	0.17	149,149,149,149	0
3	PO4	C	303	5/5	0.77	0.26	154,155,156,157	0
3	PO4	D	302	5/5	0.77	0.19	153,153,153,153	0
3	PO4	A	202	5/5	0.80	0.19	165,166,167,167	0
3	PO4	C	301	5/5	0.80	0.19	132,133,134,134	0
3	PO4	D	301	5/5	0.85	0.12	117,118,118,119	0
3	PO4	B	201	5/5	0.87	0.17	123,123,124,124	0
3	PO4	A	201	5/5	0.92	0.08	78,79,80,82	0
4	EDO	A	203	4/4	0.95	0.12	69,69,70,72	0

6.5 Other polymers [i](#)

There are no such residues in this entry.