



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 02:45 am GMT

PDB ID : 4KRY
Title : Structure of Aes from E. coli in covalent complex with PMS
Authors : Schiefner, A.; Gerber, K.; Brosig, A.; Boos, W.
Deposited on : 2013-05-17
Resolution : 2.30 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

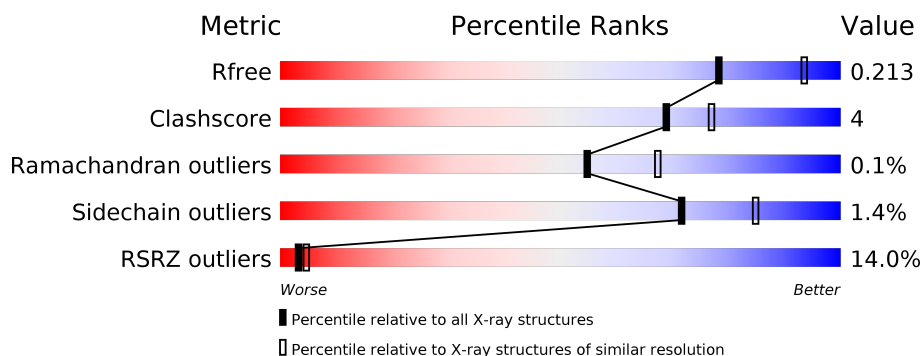
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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 $\leq 5\%$. 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	333	<div> <div>12%</div> <div>87% 9% 5%</div> </div>
1	B	333	<div> <div>10%</div> <div>87% 8% 5%</div> </div>
1	C	333	<div> <div>18%</div> <div>84% 11% 5%</div> </div>
1	D	333	<div> <div>9%</div> <div>85% 10% 5%</div> </div>
1	E	333	<div> <div>9%</div> <div>88% 11% .</div> </div>
1	F	333	<div> <div>22%</div> <div>89% 7% 5%</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	PGE	F	401	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15961 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 Acetyl esterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	0	0
			2537	1621	424	474	18			
1	B	318	Total	C	N	O	S	0	1	0
			2545	1626	427	474	18			
1	C	318	Total	C	N	O	S	0	2	0
			2554	1631	428	477	18			
1	D	318	Total	C	N	O	S	0	1	0
			2545	1626	427	474	18			
1	E	329	Total	C	N	O	S	0	3	0
			2660	1695	455	490	20			
1	F	318	Total	C	N	O	S	0	0	0
			2537	1621	424	474	18			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	EXPRESSION TAG	UNP P23872
A	-12	ARG	-	EXPRESSION TAG	UNP P23872
A	-11	GLY	-	EXPRESSION TAG	UNP P23872
A	-10	SER	-	EXPRESSION TAG	UNP P23872
A	-9	HIS	-	EXPRESSION TAG	UNP P23872
A	-8	HIS	-	EXPRESSION TAG	UNP P23872
A	-7	HIS	-	EXPRESSION TAG	UNP P23872
A	-6	HIS	-	EXPRESSION TAG	UNP P23872
A	-5	HIS	-	EXPRESSION TAG	UNP P23872
A	-4	HIS	-	EXPRESSION TAG	UNP P23872
A	-3	THR	-	EXPRESSION TAG	UNP P23872
A	-2	ASP	-	EXPRESSION TAG	UNP P23872
A	-1	PRO	-	EXPRESSION TAG	UNP P23872
A	0	ILE	-	EXPRESSION TAG	UNP P23872
B	-13	MET	-	EXPRESSION TAG	UNP P23872
B	-12	ARG	-	EXPRESSION TAG	UNP P23872
B	-11	GLY	-	EXPRESSION TAG	UNP P23872

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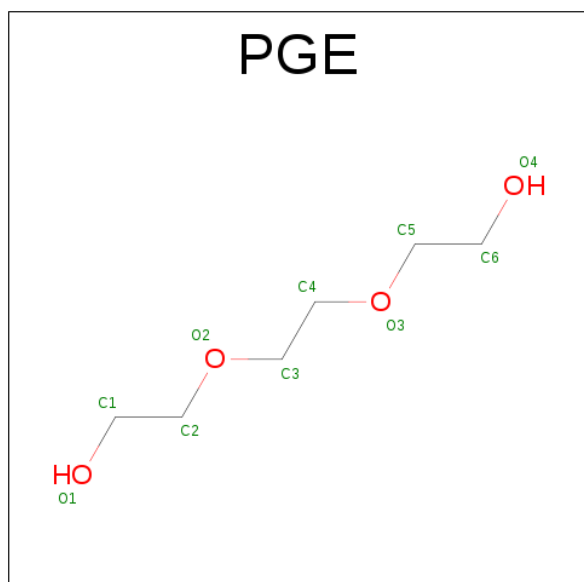
Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	SER	-	EXPRESSION TAG	UNP P23872
B	-9	HIS	-	EXPRESSION TAG	UNP P23872
B	-8	HIS	-	EXPRESSION TAG	UNP P23872
B	-7	HIS	-	EXPRESSION TAG	UNP P23872
B	-6	HIS	-	EXPRESSION TAG	UNP P23872
B	-5	HIS	-	EXPRESSION TAG	UNP P23872
B	-4	HIS	-	EXPRESSION TAG	UNP P23872
B	-3	THR	-	EXPRESSION TAG	UNP P23872
B	-2	ASP	-	EXPRESSION TAG	UNP P23872
B	-1	PRO	-	EXPRESSION TAG	UNP P23872
B	0	ILE	-	EXPRESSION TAG	UNP P23872
C	-13	MET	-	EXPRESSION TAG	UNP P23872
C	-12	ARG	-	EXPRESSION TAG	UNP P23872
C	-11	GLY	-	EXPRESSION TAG	UNP P23872
C	-10	SER	-	EXPRESSION TAG	UNP P23872
C	-9	HIS	-	EXPRESSION TAG	UNP P23872
C	-8	HIS	-	EXPRESSION TAG	UNP P23872
C	-7	HIS	-	EXPRESSION TAG	UNP P23872
C	-6	HIS	-	EXPRESSION TAG	UNP P23872
C	-5	HIS	-	EXPRESSION TAG	UNP P23872
C	-4	HIS	-	EXPRESSION TAG	UNP P23872
C	-3	THR	-	EXPRESSION TAG	UNP P23872
C	-2	ASP	-	EXPRESSION TAG	UNP P23872
C	-1	PRO	-	EXPRESSION TAG	UNP P23872
C	0	ILE	-	EXPRESSION TAG	UNP P23872
D	-13	MET	-	EXPRESSION TAG	UNP P23872
D	-12	ARG	-	EXPRESSION TAG	UNP P23872
D	-11	GLY	-	EXPRESSION TAG	UNP P23872
D	-10	SER	-	EXPRESSION TAG	UNP P23872
D	-9	HIS	-	EXPRESSION TAG	UNP P23872
D	-8	HIS	-	EXPRESSION TAG	UNP P23872
D	-7	HIS	-	EXPRESSION TAG	UNP P23872
D	-6	HIS	-	EXPRESSION TAG	UNP P23872
D	-5	HIS	-	EXPRESSION TAG	UNP P23872
D	-4	HIS	-	EXPRESSION TAG	UNP P23872
D	-3	THR	-	EXPRESSION TAG	UNP P23872
D	-2	ASP	-	EXPRESSION TAG	UNP P23872
D	-1	PRO	-	EXPRESSION TAG	UNP P23872
D	0	ILE	-	EXPRESSION TAG	UNP P23872
E	-13	MET	-	EXPRESSION TAG	UNP P23872
E	-12	ARG	-	EXPRESSION TAG	UNP P23872
E	-11	GLY	-	EXPRESSION TAG	UNP P23872

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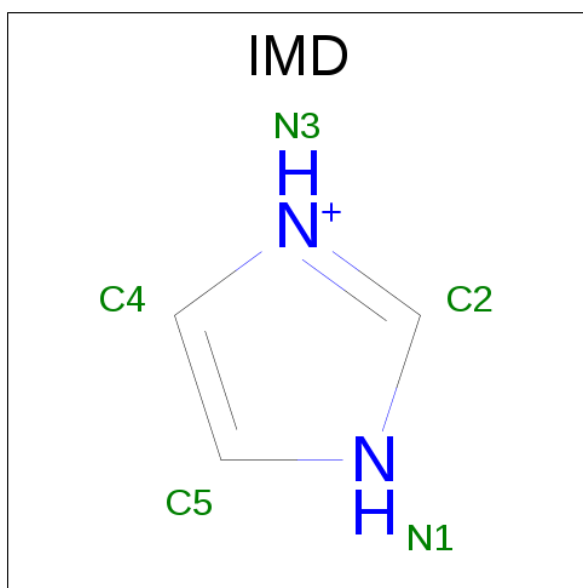
Chain	Residue	Modelled	Actual	Comment	Reference
E	-10	SER	-	EXPRESSION TAG	UNP P23872
E	-9	HIS	-	EXPRESSION TAG	UNP P23872
E	-8	HIS	-	EXPRESSION TAG	UNP P23872
E	-7	HIS	-	EXPRESSION TAG	UNP P23872
E	-6	HIS	-	EXPRESSION TAG	UNP P23872
E	-5	HIS	-	EXPRESSION TAG	UNP P23872
E	-4	HIS	-	EXPRESSION TAG	UNP P23872
E	-3	THR	-	EXPRESSION TAG	UNP P23872
E	-2	ASP	-	EXPRESSION TAG	UNP P23872
E	-1	PRO	-	EXPRESSION TAG	UNP P23872
E	0	ILE	-	EXPRESSION TAG	UNP P23872
F	-13	MET	-	EXPRESSION TAG	UNP P23872
F	-12	ARG	-	EXPRESSION TAG	UNP P23872
F	-11	GLY	-	EXPRESSION TAG	UNP P23872
F	-10	SER	-	EXPRESSION TAG	UNP P23872
F	-9	HIS	-	EXPRESSION TAG	UNP P23872
F	-8	HIS	-	EXPRESSION TAG	UNP P23872
F	-7	HIS	-	EXPRESSION TAG	UNP P23872
F	-6	HIS	-	EXPRESSION TAG	UNP P23872
F	-5	HIS	-	EXPRESSION TAG	UNP P23872
F	-4	HIS	-	EXPRESSION TAG	UNP P23872
F	-3	THR	-	EXPRESSION TAG	UNP P23872
F	-2	ASP	-	EXPRESSION TAG	UNP P23872
F	-1	PRO	-	EXPRESSION TAG	UNP P23872
F	0	ILE	-	EXPRESSION TAG	UNP P23872

- Molecule 2 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



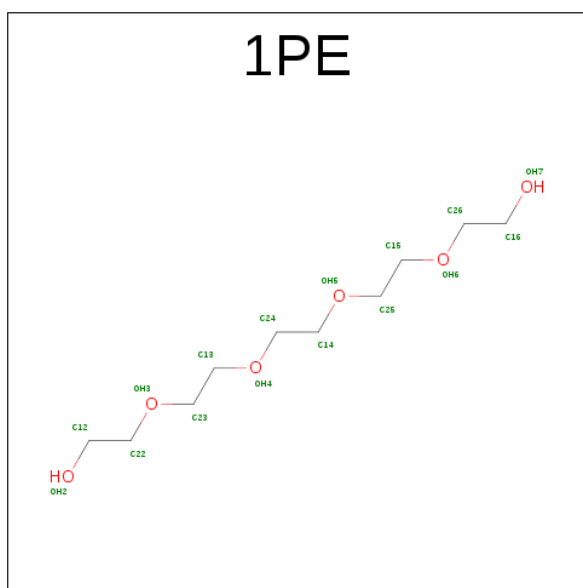
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 10 6 4	0	0
2	A	1	Total C O 10 6 4	0	0
2	B	1	Total C O 10 6 4	0	0
2	C	1	Total C O 10 6 4	0	0
2	D	1	Total C O 10 6 4	0	0
2	D	1	Total C O 10 6 4	0	0
2	E	1	Total C O 10 6 4	0	0
2	F	1	Total C O 10 6 4	0	0

- Molecule 3 is IMIDAZOLE (three-letter code: IMD) (formula: $C_3H_5N_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N 5 3 2	0	0
3	B	1	Total C N 5 3 2	0	0
3	D	1	Total C N 5 3 2	0	0
3	D	1	Total C N 5 3 2	0	0

- Molecule 4 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			16	10	6		
4	E	1	Total	C	O	0	0
			16	10	6		

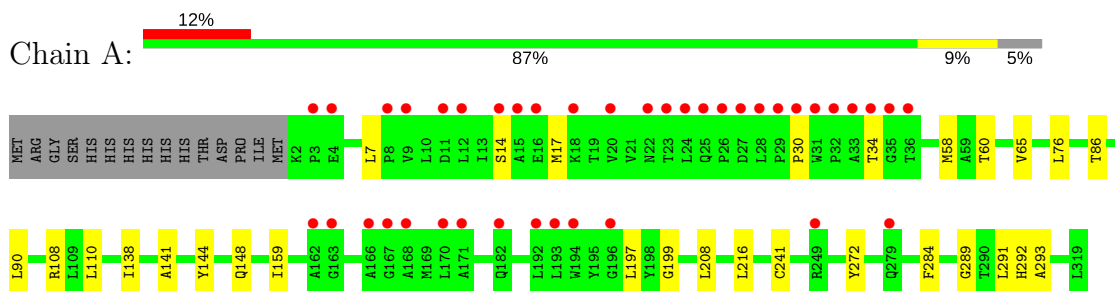
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	72	Total	O	0	0
			72	72		
5	B	82	Total	O	0	1
			83	83		
5	C	45	Total	O	0	0
			45	45		
5	D	98	Total	O	0	0
			98	98		
5	E	116	Total	O	0	1
			117	117		
5	F	36	Total	O	0	0
			36	36		

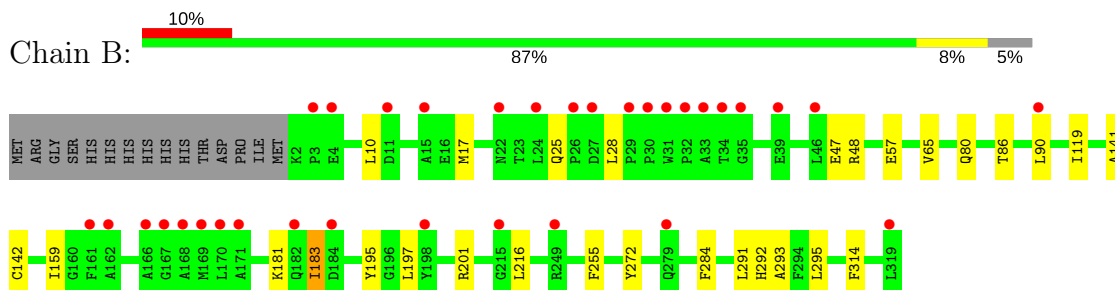
3 Residue-property plots [i](#)

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

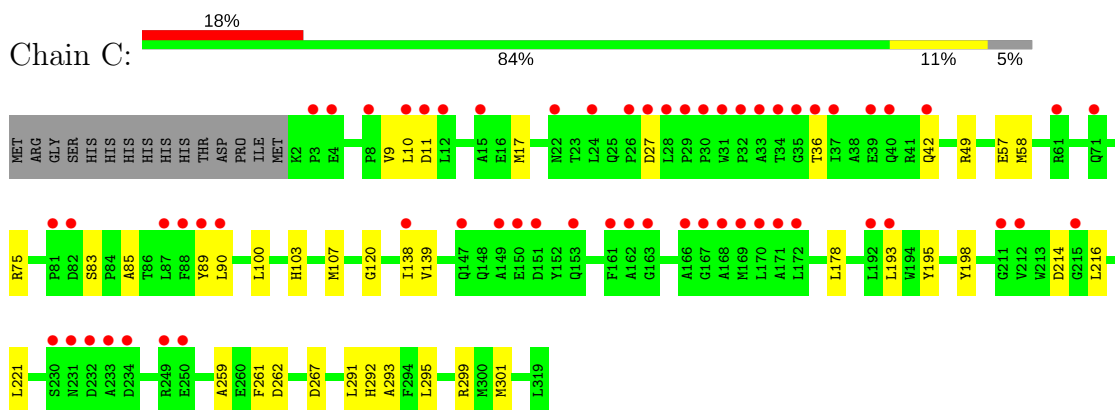
• Molecule 1: Acetyl esterase



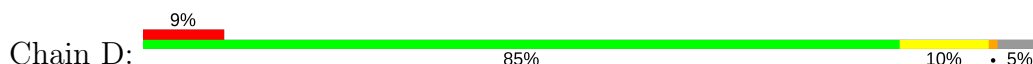
• Molecule 1: Acetyl esterase

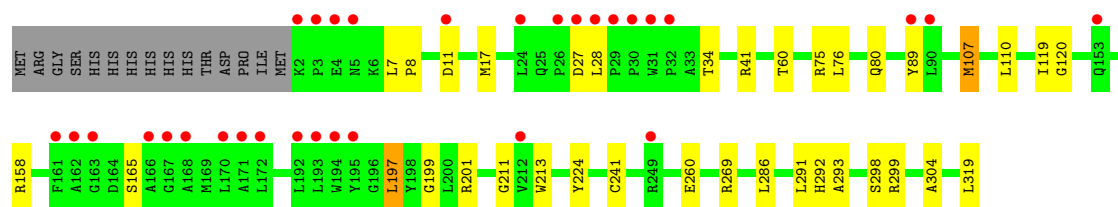


• Molecule 1: Acetyl esterase

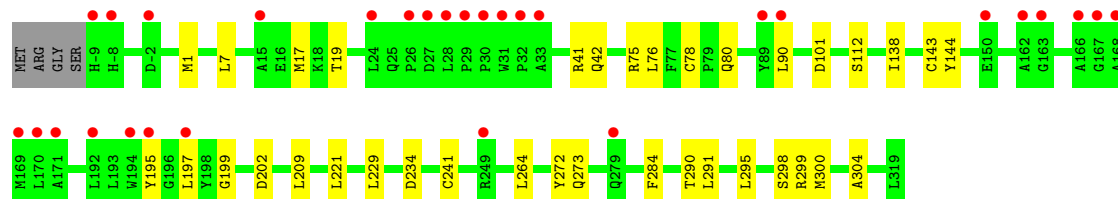
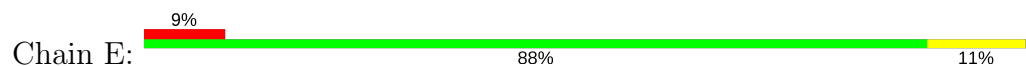


• Molecule 1: Acetyl esterase

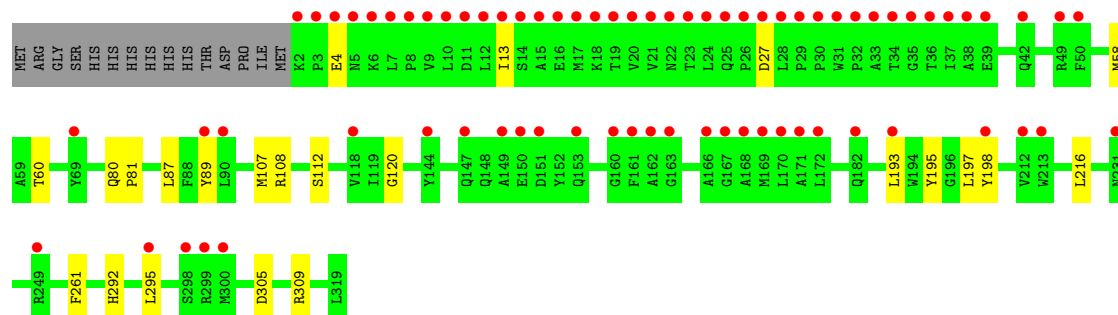
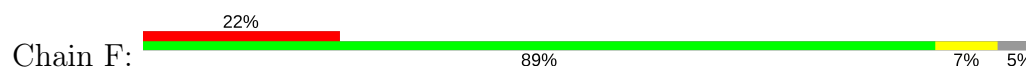




• Molecule 1: Acetyl esterase



• Molecule 1: Acetyl esterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	111.40Å 111.40Å 282.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.37 – 2.30 29.35 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.37-2.30) 99.8 (29.35-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.40 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.179 , 0.209 0.184 , 0.213	Depositor DCC
R_{free} test set	7572 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 59.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.068 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15961	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.61% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PGE, 1PE, SEB, IMD

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	0.57	0/2586	0.70	0/3516
1	B	0.61	0/2597	0.76	1/3530 (0.0%)
1	C	0.49	0/2606	0.64	0/3542
1	D	0.68	0/2597	0.81	4/3530 (0.1%)
1	E	0.64	0/2719	0.79	5/3697 (0.1%)
1	F	0.50	0/2586	0.67	1/3516 (0.0%)
All	All	0.59	0/15691	0.73	11/21331 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	41	ARG	NE-CZ-NH1	6.45	123.52	120.30
1	E	101	ASP	CB-CG-OD1	6.16	123.84	118.30
1	E	41	ARG	NE-CZ-NH2	-5.65	117.47	120.30
1	E	202	ASP	CB-CG-OD1	5.46	123.22	118.30
1	D	269	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	E	75	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	D	41	ARG	NE-CZ-NH1	5.31	122.96	120.30
1	D	201	ARG	NE-CZ-NH2	-5.25	117.67	120.30
1	B	201	ARG	NE-CZ-NH2	-5.25	117.67	120.30
1	F	87	LEU	CA-CB-CG	5.02	126.85	115.30
1	D	107	MET	CG-SD-CE	-5.01	92.18	100.20

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	2537	0	2450	19	0
1	B	2545	0	2463	20	0
1	C	2554	0	2468	26	0
1	D	2545	0	2463	23	0
1	E	2660	0	2563	25	0
1	F	2537	0	2450	19	0
2	A	20	0	28	1	0
2	B	10	0	14	1	0
2	C	10	0	14	2	0
2	D	20	0	28	3	0
2	E	10	0	14	1	0
2	F	10	0	14	2	0
3	A	5	0	5	0	0
3	B	5	0	5	0	0
3	D	10	0	10	0	0
4	D	16	0	22	0	0
4	E	16	0	22	2	0
5	A	72	0	0	0	0
5	B	83	0	0	0	0
5	C	45	0	0	0	0
5	D	98	0	0	3	0
5	E	117	0	0	2	0
5	F	36	0	0	0	0
All	All	15961	0	15033	131	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (131) 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:E:76:LEU:CD2	1:E:78[A]:CYS:SG	2.68	0.81
1:B:181:LYS:HE3	1:B:183:ILE:HD11	1.73	0.70
1:E:299[A]:ARG:HH11	1:E:299[A]:ARG:HG2	1.57	0.68
1:C:17:MET:HE3	1:C:291:LEU:H	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:197:LEU:HD12	2:F:401:PGE:H12	1.76	0.67
1:E:299[A]:ARG:NH1	1:E:299[A]:ARG:HG2	2.09	0.67
1:B:181:LYS:CE	1:B:183:ILE:HD11	2.26	0.65
1:A:197:LEU:H	1:A:197:LEU:HD23	1.62	0.64
1:B:197:LEU:HD12	2:B:401:PGE:H22	1.79	0.64
1:B:47:GLU:OE1	1:B:48:ARG:HD3	1.98	0.63
1:F:107:MET:CE	1:F:120:GLY:HA3	2.29	0.62
1:F:107:MET:HE3	1:F:120:GLY:CA	2.31	0.61
1:B:197:LEU:H	1:B:197:LEU:HD23	1.66	0.60
1:F:305:ASP:OD2	1:F:309:ARG:NH1	2.34	0.60
1:E:221:LEU:HB3	2:E:401:PGE:H32	1.83	0.59
1:B:65:VAL:HG11	1:B:141:ALA:HA	1.85	0.59
1:F:107:MET:CE	1:F:120:GLY:CA	2.81	0.59
1:B:195:TYR:OH	1:B:295:LEU:HB2	2.03	0.58
1:D:107:MET:CE	1:D:120:GLY:CA	2.80	0.58
1:E:299[A]:ARG:HH11	1:E:299[A]:ARG:CG	2.16	0.58
1:D:17:MET:HE3	1:D:291:LEU:H	1.70	0.57
1:D:197:LEU:HD12	2:D:401:PGE:H12	1.87	0.56
1:D:76:LEU:HG	1:D:119:ILE:HG12	1.87	0.56
1:A:197:LEU:HD12	2:A:401:PGE:H22	1.88	0.55
1:D:107:MET:HE3	1:D:120:GLY:CA	2.36	0.55
1:E:76:LEU:HD23	1:E:78[A]:CYS:SG	2.47	0.55
1:A:90:LEU:HD22	1:A:138:ILE:CD1	2.37	0.55
1:C:221:LEU:HB3	2:C:401:PGE:C4	2.37	0.55
1:C:139:VAL:HG13	1:C:178:LEU:HD21	1.89	0.55
2:D:402:PGE:H12	5:D:566:HOH:O	2.07	0.54
1:F:197:LEU:HD12	2:F:401:PGE:C1	2.36	0.54
1:A:144:TYR:O	1:A:148:GLN:HG2	2.08	0.54
1:C:259:ALA:HB3	1:C:262:ASP:HB2	1.91	0.53
1:F:60:THR:HG21	1:F:108:ARG:NH2	2.24	0.53
1:B:17:MET:HE3	1:B:291:LEU:H	1.73	0.53
1:C:57:GLU:O	1:C:58:MET:HB3	2.09	0.52
1:E:229:LEU:HD22	1:E:234:ASP:HB3	1.92	0.52
1:D:107:MET:CE	1:D:120:GLY:HA3	2.40	0.52
1:A:58:MET:O	1:A:60:THR:HG23	2.10	0.52
1:D:211:GLY:HA3	1:D:213:TRP:CZ3	2.45	0.52
1:A:65:VAL:HG11	1:A:141:ALA:HA	1.93	0.51
1:D:89:TYR:HB3	1:D:107:MET:HE3	1.92	0.51
1:E:90:LEU:HD22	1:E:138:ILE:CD1	2.41	0.50
1:F:107:MET:HE3	1:F:120:GLY:HA3	1.92	0.50
1:F:107:MET:HE1	1:F:120:GLY:HA3	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:SER:OG	1:C:85:ALA:O	2.30	0.49
1:D:165:SEB:HI2	1:D:224:TYR:CZ	2.48	0.49
1:E:195:TYR:OH	1:E:295:LEU:HB2	2.13	0.49
1:C:221:LEU:HB3	2:C:401:PGE:H42	1.94	0.49
1:E:199:GLY:HA3	1:E:241:CYS:HA	1.95	0.48
1:C:17:MET:HE1	1:C:301:MET:SD	2.53	0.48
1:C:90:LEU:HD22	1:C:138:ILE:CD1	2.44	0.48
1:D:107:MET:HE1	1:D:120:GLY:CA	2.44	0.48
1:A:216:LEU:HD11	1:A:292:HIS:CD2	2.48	0.48
1:A:292:HIS:O	1:A:293:ALA:HB3	2.14	0.47
1:C:89:TYR:HB3	1:C:107:MET:HE3	1.96	0.47
1:B:57:GLU:HG3	1:C:57:GLU:HG3	1.96	0.47
1:B:216:LEU:HD11	1:B:292:HIS:CG	2.49	0.47
1:D:298:SER:HA	1:D:304:ALA:HB3	1.96	0.47
2:D:402:PGE:C1	5:D:566:HOH:O	2.62	0.47
1:D:60:THR:HA	1:D:76:LEU:O	2.15	0.47
1:E:7:LEU:HD12	1:E:209:LEU:HD21	1.97	0.47
1:E:76:LEU:HD22	1:E:78[A]:CYS:SG	2.53	0.46
1:D:197:LEU:H	1:D:197:LEU:HD23	1.80	0.46
1:E:197:LEU:HD23	1:E:197:LEU:H	1.80	0.46
1:F:107:MET:HE3	1:F:120:GLY:HA2	1.96	0.46
1:E:273:GLN:NE2	5:E:596:HOH:O	2.49	0.46
1:E:298:SER:HA	1:E:304:ALA:HB3	1.98	0.46
1:C:107:MET:HE3	1:C:120:GLY:CA	2.46	0.46
1:E:143:CYS:HB3	4:E:402:1PE:H221	1.98	0.46
1:A:34:THR:O	1:A:34:THR:HG22	2.16	0.46
1:E:17:MET:HE3	1:E:291:LEU:H	1.81	0.46
1:A:110:LEU:HD12	1:A:110:LEU:HA	1.85	0.45
1:E:144:TYR:HB2	4:E:402:1PE:H131	1.98	0.45
1:B:25:GLN:HB2	1:B:28:LEU:HD12	1.97	0.45
1:D:260:GLU:HA	1:D:286:LEU:HD11	1.98	0.45
1:F:13:ILE:HG22	1:F:261:PHE:CD2	2.51	0.45
1:B:255:PHE:HB2	1:B:314:PHE:CD1	2.51	0.45
1:C:107:MET:CE	1:C:120:GLY:CA	2.95	0.45
1:D:107:MET:HE1	1:D:120:GLY:HA3	1.97	0.45
1:B:216:LEU:HD11	1:B:292:HIS:CD2	2.52	0.45
1:D:107:MET:HE1	1:D:120:GLY:N	2.31	0.45
1:C:9:VAL:HG21	1:C:261:PHE:HB3	1.99	0.45
1:F:89:TYR:HB3	1:F:107:MET:CE	2.47	0.45
1:F:197:LEU:HD23	1:F:197:LEU:H	1.81	0.45
1:F:58:MET:O	1:F:60:THR:HG23	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:195:TYR:OH	1:F:295:LEU:HB2	2.16	0.45
1:C:103:HIS:O	1:C:107:MET:HG3	2.18	0.44
1:D:199:GLY:HA3	1:D:241:CYS:HA	1.99	0.44
1:C:299[B]:ARG:HH11	1:C:299[B]:ARG:HG2	1.83	0.44
1:A:272:TYR:CD1	1:A:284:PHE:HB2	2.52	0.44
1:F:80:GLN:HB2	1:F:81:PRO:HD2	2.00	0.44
1:A:17:MET:HE3	1:A:291:LEU:H	1.83	0.43
1:C:299[B]:ARG:HG2	1:C:299[B]:ARG:NH1	2.33	0.43
1:D:75:ARG:NH1	5:D:536:HOH:O	2.47	0.43
1:C:89:TYR:HB3	1:C:107:MET:CE	2.48	0.43
1:D:299[B]:ARG:HD3	1:E:112:SER:O	2.18	0.43
1:E:272:TYR:CD1	1:E:284:PHE:HB2	2.54	0.43
1:A:86:THR:O	1:A:159:ILE:HA	2.18	0.43
1:A:199:GLY:HA3	1:A:241:CYS:HA	1.99	0.43
1:A:60:THR:HG21	1:A:108:ARG:NH2	2.33	0.43
1:F:107:MET:HE1	1:F:120:GLY:CA	2.47	0.43
1:A:216:LEU:HD11	1:A:292:HIS:CG	2.54	0.42
1:B:86:THR:O	1:B:159:ILE:HA	2.18	0.42
1:C:195:TYR:OH	1:C:295:LEU:HB2	2.19	0.42
1:A:60:THR:HA	1:A:76:LEU:O	2.20	0.42
1:E:42:GLN:NE2	5:E:578:HOH:O	2.53	0.42
1:C:216:LEU:HD11	1:C:292:HIS:CD2	2.54	0.42
1:D:158:ARG:NH1	1:D:319:LEU:O	2.53	0.42
1:E:17:MET:HE2	1:E:290:THR:HA	2.00	0.42
1:C:292:HIS:O	1:C:293:ALA:HB3	2.20	0.42
1:B:90:LEU:HD22	1:B:90:LEU:N	2.35	0.41
1:D:7:LEU:HD12	1:D:8:PRO:HD2	2.01	0.41
1:D:292:HIS:O	1:D:293:ALA:HB3	2.21	0.41
1:F:216:LEU:HD11	1:F:292:HIS:CG	2.56	0.41
1:C:75:ARG:HB2	1:C:100:LEU:HD13	2.03	0.41
1:B:272:TYR:CD1	1:B:284:PHE:HB2	2.56	0.41
1:B:292:HIS:O	1:B:293:ALA:HB3	2.21	0.41
1:B:86:THR:HG21	1:B:119:ILE:HD12	2.02	0.41
1:E:17:MET:CE	1:E:290:THR:HA	2.51	0.41
1:A:14:SER:HB3	1:A:289:GLY:HA3	2.03	0.41
1:C:107:MET:HE3	1:C:120:GLY:HA2	2.03	0.41
1:D:107:MET:CE	1:D:120:GLY:N	2.84	0.41
1:E:19:THR:HG22	1:E:300:MET:CE	2.52	0.40
1:B:181:LYS:HE2	1:B:183:ILE:HD11	2.02	0.40
1:A:7:LEU:HD22	1:A:208:LEU:HG	2.02	0.40
1:E:197:LEU:HB3	1:E:264:LEU:HD13	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:193:LEU:HB3	1:F:198:TYR:OH	2.22	0.40
1:B:57:GLU:HG3	1:C:57:GLU:CG	2.52	0.40
1:C:107:MET:CE	1:C:120:GLY:HA3	2.52	0.40
1:C:193:LEU:HB3	1:C:198:TYR:OH	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	315/333 (95%)	306 (97%)	9 (3%)	0	100	100
1	B	316/333 (95%)	302 (96%)	14 (4%)	0	100	100
1	C	317/333 (95%)	302 (95%)	15 (5%)	0	100	100
1	D	316/333 (95%)	302 (96%)	14 (4%)	0	100	100
1	E	329/333 (99%)	317 (96%)	11 (3%)	1 (0%)	44	55
1	F	315/333 (95%)	303 (96%)	11 (4%)	1 (0%)	44	55
All	All	1908/1998 (96%)	1832 (96%)	74 (4%)	2 (0%)	55	67

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	1	MET
1	F	4	GLU

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	264/278 (95%)	263 (100%)	1 (0%)	93	97
1	B	265/278 (95%)	261 (98%)	4 (2%)	70	83
1	C	266/278 (96%)	258 (97%)	8 (3%)	46	63
1	D	265/278 (95%)	258 (97%)	7 (3%)	51	69
1	E	278/278 (100%)	277 (100%)	1 (0%)	93	97
1	F	264/278 (95%)	262 (99%)	2 (1%)	85	93
All	All	1602/1668 (96%)	1579 (99%)	23 (1%)	71	85

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	PRO
1	B	10	LEU
1	B	80	GLN
1	B	142	CYS
1	B	183	ILE
1	C	10	LEU
1	C	11	ASP
1	C	27	ASP
1	C	36	THR
1	C	42	GLN
1	C	49	ARG
1	C	214	ASP
1	C	267	ASP
1	D	11	ASP
1	D	27	ASP
1	D	28	LEU
1	D	34	THR
1	D	80	GLN
1	D	110	LEU
1	D	197	LEU
1	E	80	GLN
1	F	27	ASP
1	F	112	SER

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	SEB	A	165	1	15,16,17	1.70	1 (6%)	16,21,23	1.16	1 (6%)
1	SEB	B	165	1	15,16,17	1.20	1 (6%)	16,21,23	0.87	1 (6%)
1	SEB	C	165	1	15,16,17	1.34	1 (6%)	16,21,23	1.04	2 (12%)
1	SEB	D	165	1	15,16,17	1.12	1 (6%)	16,21,23	1.06	2 (12%)
1	SEB	E	165	1	15,16,17	0.92	1 (6%)	16,21,23	0.91	1 (6%)
1	SEB	F	165	1	15,16,17	1.16	1 (6%)	16,21,23	1.05	1 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	SEB	A	165	1	-	0/9/13/15	0/1/1/1
1	SEB	B	165	1	-	0/9/13/15	0/1/1/1
1	SEB	C	165	1	-	0/9/13/15	0/1/1/1
1	SEB	D	165	1	-	0/9/13/15	0/1/1/1
1	SEB	E	165	1	-	0/9/13/15	0/1/1/1
1	SEB	F	165	1	-	0/9/13/15	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	165	SEB	CA-C	2.24	1.53	1.50
1	D	165	SEB	CA-C	3.49	1.54	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	165	SEB	CA-C	3.66	1.55	1.50
1	B	165	SEB	CA-C	4.10	1.55	1.50
1	C	165	SEB	CA-C	4.63	1.56	1.50
1	A	165	SEB	CA-C	5.88	1.58	1.50

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	165	SEB	O-C-CA	-2.23	118.85	125.02
1	D	165	SEB	O-C-CA	-2.18	118.99	125.02
1	B	165	SEB	O-C-CA	-2.14	119.10	125.02
1	C	165	SEB	O-C-CA	-2.12	119.17	125.02
1	C	165	SEB	CB-OG-SD	2.02	123.33	119.12
1	D	165	SEB	CB-OG-SD	2.24	123.78	119.12
1	F	165	SEB	OG-SD-CE	2.69	111.54	104.08
1	A	165	SEB	OD1-SD-CE	2.76	114.82	108.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	165	SEB	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

14 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
2	PGE	A	401	-	9,9,9	0.54	0	8,8,8	0.28	0
2	PGE	A	402	-	9,9,9	0.75	0	8,8,8	0.59	0
3	IMD	A	403	-	3,5,5	0.41	0	4,5,5	0.65	0
2	PGE	B	401	-	9,9,9	0.51	0	8,8,8	0.42	0
3	IMD	B	402	-	3,5,5	0.45	0	4,5,5	0.60	0
2	PGE	C	401	-	9,9,9	0.39	0	8,8,8	0.28	0
2	PGE	D	401	-	9,9,9	0.71	0	8,8,8	0.59	0
2	PGE	D	402	-	9,9,9	0.56	0	8,8,8	0.47	0
4	1PE	D	403	-	15,15,15	0.57	0	14,14,14	0.41	0
3	IMD	D	404	-	3,5,5	0.43	0	4,5,5	0.64	0
3	IMD	D	405	-	3,5,5	0.59	0	4,5,5	0.38	0
2	PGE	E	401	-	9,9,9	0.64	0	8,8,8	0.53	0
4	1PE	E	402	-	15,15,15	0.75	0	14,14,14	0.90	0
2	PGE	F	401	-	9,9,9	0.49	0	8,8,8	0.83	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
2	PGE	A	401	-	-	0/7/7/7	0/0/0/0
2	PGE	A	402	-	-	0/7/7/7	0/0/0/0
3	IMD	A	403	-	-	0/0/0/0	0/1/1/1
2	PGE	B	401	-	-	0/7/7/7	0/0/0/0
3	IMD	B	402	-	-	0/0/0/0	0/1/1/1
2	PGE	C	401	-	-	0/7/7/7	0/0/0/0
2	PGE	D	401	-	-	0/7/7/7	0/0/0/0
2	PGE	D	402	-	-	0/7/7/7	0/0/0/0
4	1PE	D	403	-	-	0/13/13/13	0/0/0/0
3	IMD	D	404	-	-	0/0/0/0	0/1/1/1
3	IMD	D	405	-	-	0/0/0/0	0/1/1/1
2	PGE	E	401	-	-	0/7/7/7	0/0/0/0
4	1PE	E	402	-	-	0/13/13/13	0/0/0/0
2	PGE	F	401	-	-	0/7/7/7	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.

8 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	PGE	1	0
2	B	401	PGE	1	0
2	C	401	PGE	2	0
2	D	401	PGE	1	0
2	D	402	PGE	2	0
2	E	401	PGE	1	0
4	E	402	1PE	2	0
2	F	401	PGE	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

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, 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	317/333 (95%)	0.45	40 (12%) 4 6	25, 39, 88, 113	0
1	B	317/333 (95%)	0.33	33 (10%) 7 10	27, 37, 74, 115	0
1	C	317/333 (95%)	0.84	60 (18%) 1 2	32, 53, 95, 134	0
1	D	317/333 (95%)	0.37	31 (9%) 8 11	24, 31, 83, 128	0
1	E	328/333 (98%)	0.25	30 (9%) 10 14	24, 32, 52, 107	0
1	F	317/333 (95%)	1.05	73 (23%) 1 1	33, 50, 122, 144	0
All	All	1913/1998 (95%)	0.55	267 (13%) 3 5	24, 39, 91, 144	0

All (267) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	3	PRO	11.0
1	F	3	PRO	9.2
1	C	3	PRO	8.8
1	F	30	PRO	8.8
1	C	34	THR	8.6
1	B	33	ALA	8.5
1	C	29	PRO	8.3
1	F	34	THR	8.0
1	C	33	ALA	7.9
1	F	28	LEU	7.9
1	F	12	LEU	7.7
1	D	26	PRO	7.5
1	F	27	ASP	7.3
1	F	26	PRO	7.2
1	C	30	PRO	7.1
1	F	33	ALA	7.0
1	F	11	ASP	6.9
1	E	26	PRO	6.6
1	F	17	MET	6.5

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Mol	Chain	Res	Type	RSRZ
1	C	26	PRO	6.4
1	B	34	THR	6.3
1	F	15	ALA	6.2
1	A	34	THR	6.2
1	F	21	VAL	6.2
1	F	31	TRP	6.1
1	F	7	LEU	6.0
1	A	33	ALA	5.9
1	F	32	PRO	5.9
1	C	28	LEU	5.7
1	E	-9	HIS	5.7
1	F	2	LYS	5.7
1	B	26	PRO	5.6
1	C	32	PRO	5.6
1	F	9	VAL	5.5
1	F	8	PRO	5.5
1	D	30	PRO	5.5
1	A	24	LEU	5.4
1	C	24	LEU	5.3
1	F	29	PRO	5.2
1	A	30	PRO	5.2
1	B	30	PRO	5.2
1	F	212	VAL	5.1
1	C	171	ALA	5.0
1	D	32	PRO	4.9
1	F	20	VAL	4.9
1	E	31	TRP	4.9
1	F	19	THR	4.8
1	D	212	VAL	4.8
1	F	24	LEU	4.8
1	C	31	TRP	4.8
1	C	150	GLU	4.7
1	A	26	PRO	4.7
1	B	31	TRP	4.7
1	B	32	PRO	4.7
1	C	27	ASP	4.7
1	D	4	GLU	4.7
1	A	32	PRO	4.6
1	C	90	LEU	4.5
1	D	2	LYS	4.5
1	D	31	TRP	4.5
1	A	31	TRP	4.5

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Mol	Chain	Res	Type	RSRZ
1	F	300	MET	4.3
1	F	37	ILE	4.3
1	A	15	ALA	4.3
1	A	27	ASP	4.2
1	C	37	ILE	4.2
1	F	25	GLN	4.1
1	F	23	THR	4.1
1	F	36	THR	4.0
1	F	4	GLU	4.0
1	F	90	LEU	4.0
1	C	215	GLY	4.0
1	E	-8	HIS	4.0
1	A	28	LEU	4.0
1	F	22	ASN	3.9
1	A	29	PRO	3.8
1	F	162	ALA	3.8
1	A	3	PRO	3.8
1	C	147	GLN	3.8
1	F	171	ALA	3.7
1	D	24	LEU	3.7
1	C	11	ASP	3.7
1	E	33	ALA	3.7
1	C	230	SER	3.7
1	D	90	LEU	3.7
1	C	170	LEU	3.7
1	C	151	ASP	3.6
1	A	12	LEU	3.6
1	A	9	VAL	3.6
1	E	24	LEU	3.5
1	F	299	ARG	3.5
1	C	10	LEU	3.5
1	D	170	LEU	3.5
1	C	8	PRO	3.5
1	C	36	THR	3.5
1	A	182	GLN	3.5
1	E	30	PRO	3.4
1	D	163	GLY	3.4
1	B	4	GLU	3.4
1	C	249	ARG	3.4
1	F	168	ALA	3.4
1	F	14	SER	3.4
1	D	5	ASN	3.4

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Mol	Chain	Res	Type	RSRZ
1	E	162	ALA	3.4
1	F	151	ASP	3.4
1	C	172	LEU	3.4
1	C	4	GLU	3.3
1	E	15	ALA	3.3
1	B	27	ASP	3.3
1	E	32	PRO	3.3
1	C	231	ASN	3.3
1	F	42	GLN	3.3
1	C	89	TYR	3.3
1	E	279	GLN	3.3
1	F	182	GLN	3.2
1	F	13	ILE	3.2
1	D	171	ALA	3.2
1	E	168	ALA	3.2
1	C	192	LEU	3.2
1	B	168	ALA	3.2
1	C	162	ALA	3.2
1	F	18	LYS	3.2
1	A	11	ASP	3.2
1	B	15	ALA	3.2
1	D	162	ALA	3.2
1	B	171	ALA	3.1
1	F	5	ASN	3.1
1	E	89	TYR	3.1
1	C	193	LEU	3.1
1	F	89	TYR	3.1
1	B	24	LEU	3.1
1	F	193	LEU	3.1
1	A	168	ALA	3.0
1	B	29	PRO	3.0
1	A	162	ALA	3.0
1	E	27	ASP	3.0
1	E	170	LEU	3.0
1	F	39	GLU	3.0
1	C	163	GLY	3.0
1	E	163	GLY	3.0
1	D	89	TYR	3.0
1	F	172	LEU	3.0
1	C	42	GLN	3.0
1	F	170	LEU	2.9
1	A	171	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	88	PHE	2.9
1	F	150	GLU	2.9
1	C	233	ALA	2.9
1	A	22	ASN	2.9
1	F	161	PHE	2.9
1	E	166	ALA	2.9
1	D	27	ASP	2.9
1	A	23	THR	2.8
1	A	36	THR	2.8
1	C	35	GLY	2.8
1	C	153	GLN	2.8
1	F	6	LYS	2.8
1	F	35	GLY	2.8
1	D	11	ASP	2.8
1	D	29	PRO	2.8
1	A	249	ARG	2.8
1	C	15	ALA	2.8
1	A	4	GLU	2.8
1	C	167	GLY	2.8
1	D	192	LEU	2.8
1	D	193	LEU	2.8
1	C	12	LEU	2.7
1	E	28	LEU	2.7
1	F	298	SER	2.7
1	F	153	GLN	2.7
1	F	38	ALA	2.7
1	C	161	PHE	2.7
1	C	39	GLU	2.6
1	B	90	LEU	2.6
1	F	231	ASN	2.6
1	A	192	LEU	2.6
1	D	168	ALA	2.6
1	C	71	GLN	2.5
1	F	147	GLN	2.5
1	C	211	GLY	2.5
1	E	90	LEU	2.5
1	A	167	GLY	2.5
1	F	167	GLY	2.5
1	F	295	LEU	2.5
1	D	249	ARG	2.5
1	F	163	GLY	2.5
1	F	198	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	162	ALA	2.4
1	E	171	ALA	2.4
1	D	194	TRP	2.4
1	B	11	ASP	2.4
1	B	279	GLN	2.4
1	C	82	ASP	2.4
1	B	39	GLU	2.4
1	E	192	LEU	2.4
1	F	50	PHE	2.4
1	A	25	GLN	2.4
1	B	182	GLN	2.4
1	D	28	LEU	2.4
1	F	166	ALA	2.4
1	A	20	VAL	2.4
1	C	212	VAL	2.4
1	E	-2	ASP	2.3
1	C	168	ALA	2.3
1	A	196	GLY	2.3
1	B	166	ALA	2.3
1	D	166	ALA	2.3
1	E	167	GLY	2.3
1	F	49	ARG	2.3
1	E	195	TYR	2.3
1	E	197	LEU	2.3
1	B	3	PRO	2.3
1	A	193	LEU	2.3
1	A	18	LYS	2.2
1	C	138	ILE	2.2
1	F	149	ALA	2.2
1	B	170	LEU	2.2
1	F	118	VAL	2.2
1	B	184	ASP	2.2
1	B	319	LEU	2.2
1	F	10	LEU	2.2
1	F	160	GLY	2.2
1	F	169	MET	2.2
1	E	29	PRO	2.2
1	C	166	ALA	2.2
1	C	232	ASP	2.2
1	D	153	GLN	2.2
1	F	249	ARG	2.2
1	C	87	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	22	ASN	2.1
1	D	167	GLY	2.2
1	C	61	ARG	2.1
1	C	250	GLU	2.1
1	B	46	LEU	2.1
1	C	169	MET	2.1
1	E	169	MET	2.1
1	E	194	TRP	2.1
1	A	166	ALA	2.1
1	F	69	TYR	2.1
1	A	14	SER	2.1
1	A	16	GLU	2.1
1	A	35	GLY	2.1
1	F	16	GLU	2.1
1	D	195	TYR	2.1
1	A	8	PRO	2.1
1	B	35	GLY	2.1
1	B	167	GLY	2.1
1	A	170	LEU	2.1
1	D	172	LEU	2.1
1	B	249	ARG	2.1
1	C	149	ALA	2.1
1	F	213	TRP	2.1
1	A	163	GLY	2.1
1	C	81	PRO	2.1
1	E	249	ARG	2.1
1	A	194	TRP	2.1
1	B	161	PHE	2.1
1	D	161	PHE	2.1
1	C	234	ASP	2.0
1	C	22	ASN	2.0
1	B	215	GLY	2.0
1	B	198	TYR	2.0
1	F	144	TYR	2.0
1	E	150	GLU	2.0
1	A	279	GLN	2.0
1	B	169	MET	2.0
1	C	40	GLN	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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, 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	LLDF	B-factors(Å ²)	Q<0.9
1	SEB	E	165	16/17	0.95	0.15	-	28,36,38,38	0
1	SEB	B	165	16/17	0.97	0.14	-	33,43,48,49	0
1	SEB	D	165	16/17	0.94	0.15	-	26,32,35,35	0
1	SEB	A	165	16/17	0.96	0.17	-	33,44,48,48	0
1	SEB	C	165	16/17	0.94	0.17	-	41,59,67,68	0
1	SEB	F	165	16/17	0.94	0.16	-	39,54,64,64	0

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, 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	LLDF	B-factors(Å ²)	Q<0.9
2	PGE	F	401	10/10	0.88	0.30	2.24	47,57,73,75	0
4	1PE	E	402	16/16	0.84	0.17	1.92	48,71,86,87	0
2	PGE	A	402	10/10	0.76	0.20	1.87	53,61,64,65	0
2	PGE	D	402	10/10	0.91	0.17	1.45	40,43,55,56	0
2	PGE	D	401	10/10	0.91	0.19	0.91	44,46,64,64	0
2	PGE	E	401	10/10	0.94	0.20	0.77	34,42,56,58	0
2	PGE	A	401	10/10	0.87	0.20	0.61	55,61,77,78	0
2	PGE	B	401	10/10	0.90	0.20	0.57	42,53,73,76	0
4	1PE	D	403	16/16	0.94	0.15	0.32	48,61,64,64	0
2	PGE	C	401	10/10	0.96	0.17	0.31	44,53,66,70	0
3	IMD	B	402	5/5	0.96	0.11	-0.72	47,49,50,52	0
3	IMD	D	405	5/5	0.97	0.10	-2.28	43,45,48,48	0
3	IMD	A	403	5/5	0.77	0.21	-	85,88,91,93	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	IMD	D	404	5/5	0.93	0.21	-	62,63,64,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.